EPA Region 5 Records Ctr.

SITE ASSESSMENT REPORT FOR STANDARD SCRAP METAL / CHICAGO INTERNATIONAL EXPORTING (aka: SCRAP METAL) U.S. EPA ID: ILD045698263 88 ID: NA TDD: T05-9402-007 PAN: EIL08318AA

MAY 6, 1994

Reviewed by:_

Approved by

PLAINTIFF'S

NOTICE

Originally, this site was tasked to Ecology & Environment, Inc. under the name Scrap Metal. The actual name of this site is Standard Scrap Metal / Chicago International Exporting. All file information, photologs, maps, and text in this report refer to the site by its original tasked name Scrap Metal.

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1.0 INTRODUCTION

The Ecology & Environment, Inc. (E & E), Technical Assistance Team (TAT) was tasked by the United States Environmental Protection Agency (U.S. EPA) to perform a site assessment at the Scrap Metal (SM) site under Technical Directive Document (TDD) T05-9402-007. Activities performed for the site assessment included prepared and implemented Health and Safety Plan; compiled available information; prepared and implemented sampling plan; evaluated threat to human health and environment; and provided photodocumentation. The following report provides a summary of these activities.

2.0 SITE DESCRIPTION

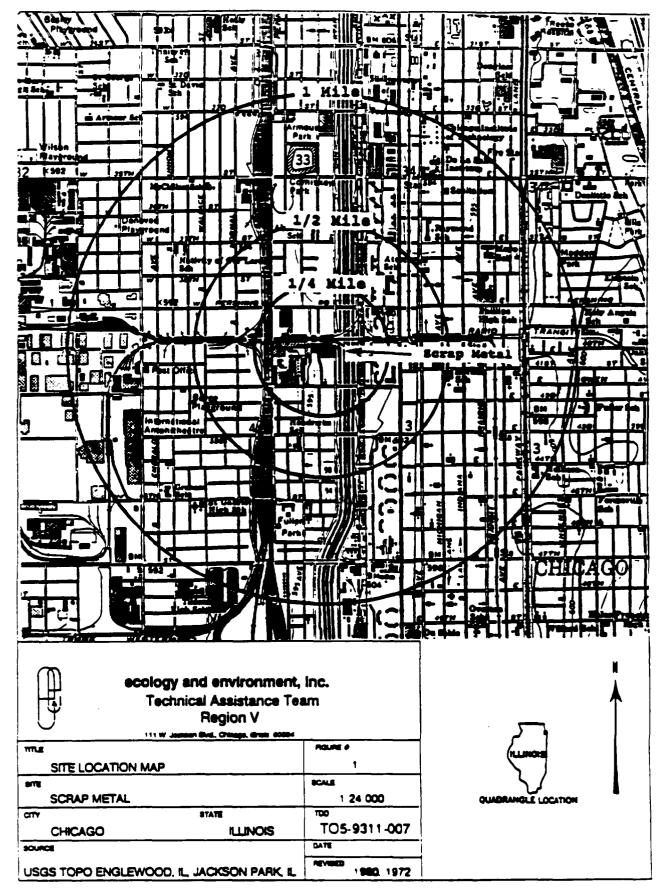
The SM site is located at 4004 South Wentworth Avenue, Chicago, Cook County, Illinois, Latitude 87° 37' 55" north, Longitude 41° 52' 50" west, (see Figure 1, Site Location Map). The SM site is located in a industrial and residential area. The facility is an active 3-acre scrap yard involved in the reclamation of metals. Past and present operations have taken place on two distinct parcels of property separated by Wells Avenue.

The east lot is approximately 2.5 acres in size and the west lot is approximately 0.5 acres (see Figure 2, Site Features map). The east lot has a office building, a small receiving shed, two aluminum furnaces, a ferrous metal sorter, a shredder, and large piles of scrap metal located at various points throughout the property. The west lot has a small scale house used to weigh incoming scrap metal trucks and is also used to park semitrailers. A railroad spur cuts across both lots diagonally.

3.0 SITE BACKGROUND

The property has been used for industrial purposes since 1895. A Sanborn Fire Insurance map, dated, 1895 indicates part of the east and west lots was owned by Weaver and Getz. It is unknown what business Weaver and Getz operated. Two other businesses owned parts of the east lot; W. B. Scace and Company (which loaded lime and cement) and Chas D. Pek and Company (which sold hay, grain, and feed). A Sanborn Fire Insurance map, dated, 1925, shows a Baker-Smith Coal Company operated a coal yard on the east lot. The west lot did not appear to be used by either Chicago Mfg. & Distribution Company (a machine shop) or Arthur M. Adler & Company, Oils located to the south.

The Standard Scrap Metal Company (SSMC) was started in 1928 by Sam Cohen and Sam Kanter at 4004 Wentworth Avenue. SSMC was involved in reclaiming aluminum and cooper and sold the reclaimed scrap metal to steel smelters and refiners. The facility contained 1 gas-fired boiler, 2 aluminum sweat furnaces, and a



wire burning incinerator. Operations continued until 1972 when the company merged into Standard Scrap Metal Company, Incorporated (SSCMI). SSCMI continued operations at the site until 1987, when the company went bankrupt.

Phoenix Recycling, a metals reclamation business, began operations at the site about the same time of the SSCMI bankruptcy and continued operating the same type of business until 1989. The Sam Cohen and Sam and Benjamin Kanter Building Partnership owned Phoenix Recycling.

Chicago International Exporting Company (CIEC) began operations at the site in 1989 and continues to operate a metal recycling facility. CIEC is owned by Chicago International Export Company, Incorporated whose President is Steve Cohen.

The SM site has been investigated by the Illinois Environmental Protection Agency (IEPA) and U.S. EPA beginning in 1973. In 1973, personnel from IEPA inspected the site for compliance with Air Pollution Regulations. The inspection revealed that the facility did not have the proper air pollution permits to operate their incinerator or sweat furnaces. A suit (PCB 83-22) was filed against SSMCI for not possessing permits required by IEPA and the City of Chicago. The complaint stated that SSMCI could achieve compliance by installing afterburners on the sweat furnaces. The afterburners were not installed and permits were not applied for until 1984. A permit for the gas-fired boiler was applied for and approved on December 14, 1984.

On January 10, 1985, Illinois Pollution Board (IPB) continued the suit (PCB 83-22) against SSMCI for permit violations. IPB suit ordered SSMCI to:

Cease and desist from operations of its incinerator until the necessary operating permit is obtained from the IEPA:

Cease and desist from operating either of its aluminum sweat furnaces until the necessary permits are obtained from the IEPA and permanently shut down the inactive aluminum sweat furnace by January 21, 1985.

Install temperature gauges on each afterburner with an interlock that prevents operation unless the afterburner temperature is at least 1400 degrees Fahrenheit, and take all necessary steps to ensure adequate pre-heating of each afterburner prior to charging. These requirements are to be made conditions of the operating permits issued by the IEPA;

Within 90 days of the date of this order pay a penalty of \$30,000 for the violation of the Act and Regulations as described in this opinion.

On February 14, 1984, IEPA investigated a report from an employee of Heatbath Corporation, the plant south of the west lot, that workers at the facility periodically dumped transformer oil on the ground and igniting it. This practice was to have taken place from 1977 to 1981. On one occasion the roof of the Heatbath Corporation caught fire and the Chicago Fire Department extinguished the fire.

IEPA collected 2 soil samples, one sample from the west lot and one from near a garage located at 3949 South Wells Avenue. A complaint from the resident at 3949 South Wells Avenue stated oil from SM Site would flow off-site into her yard. Sampling results indicated Polychlorinated biphenyls (PCBs) at 3.9 parts per million (ppm) at the residence and 1,300 ppm from the west lot. The IEPA requested that the U.S. EPA conduct a PCB inspection at the site.

On March 30, 1984, U.S. EPA's Toxic Substance Office conducted a inspection of SM facility to document their handling, storage, and disposal practices. During the inspection 6 composite soil samples and 1 wipe sample were collected from the west lot and a residence at 3949 South Wells Avenue. Analytical results from the samples indicated PCB levels of up to 2,095 ppm but no detectable amounts at the residence. The U.S. EPA filed a complaint against SSMCI for violating regulations pertaining to the disposal of PCBs and a civil penalty of \$25,000 was levied for improper disposal of PCBs.

On June 18, 1985, the Roy F. Weston TAT collected 4 soil samples and 2 wipe samples (east lot) at the SM site. The analytical results indicated PCBs levels up to 336 ppm in three samples and some dioxin isomers in 4 samples.

On October 29, 1985, an amended complaint by U.S. EPA was filed against SSMCI after the June 18, 1985 inspection. The amended complaint levied a \$30,000 penalty for violations of Section 16(a) of the Toxic Substance Control Act (TSCA). In February 1987, SSMCI appealed the decision and the complaint was dismissed because the U.S. EPA did not prove that the PCB-contaminated oil had been accepted at the site after 1978. However, the U.S. EPA appealed the dismissal and the decision was reversed and the \$30,000 fine was levied against the facility. SSMCI filed for bankruptcy, the fine was not collected and the U.S. EPA Enforcement Branch closed the case. In 1989, metal recycling operations continued on-site, when CIEC began operation at the site.

In 1990, a former railroad employee had a telephone interview with Tom Crause of the IEPA. The former railroad employee indicated that workers at the facility cut up and disposed of many electrical transformers during his 30 years of working for the railroad. Based on this information, on August 27, 1990,

SSMC was placed on the Comprehensive Environmental Response, Compensation and Liability Information System (CERCLIS).

On August 29, 1991, IEPA personnel conducted an off-site reconnaissance inspection of the SM facility. IEPA observed piles of scrap metal around the site. No air admission were observed at the site, the boiler appeared to be not in operation. At the east lot, the north sweat furnace had been demolished and was left as a pile of debris. A number of drums, which appeared to be empty, were observed near the north side of the office building. No leakage was observed from the drums or stressed vegetation on the lot. At the west lot, the gates were open and was empty except for 3 semi-trailers. The IEPA prepared a Preliminary Assessment (PA) for Standard Scrap Metal site (also called Standard Metal) on September 30, 1991.

On September 22, 1992, the IEPA was tasked by the U.S. EPA Region V to conduct a CERCLIS Screening Site Inspection (SSI) of the SM site. The SSI was conducted on November 4 and 5, 1992 and consisted of the collection of 12 soil samples. The analytical results from on-site soil sampling indicated PCBs up to 670,000 Micrograms per kilogram (ug/kg = parts per billion (ppb). The PCBs can be directly associated with past activities at the site as reported by the Heatbath employee and railroad employee.

4.0 SITE ASSESSMENT

On February 22, 1994, TAT members John Nordine and Yvette Anderson met with U.S. EPA On-Scene-Coordinator (OSC) Steve Faryan at the SM site. The SM site is currently called Chicago International Export Company. Access to the site was granted during an interview with CIEC President, Steve Cohn and Secretary Treasure, Buddy Cohn. A site safety meeting was held and the potential hazards of the site were discussed. A site reconnaissance of the SM Site was completed, which included a tour of the facility (see Appendix A for Site Photo Log).

A railroad spur cuts across both lots diagonally. A railroad overpass delineates the north property line for both lots (see Figure 2, Site Features map). A 8-foot fence topped with 3 strands of barbed wire surround both lots.

The east lot has an office building, a small receiving shed, 2 aluminum furnaces, a ferrous metal sorter, a shredder, a ash pile, and large piles of scrap metal located at various points throughout the property. According to Buddy Cohn, the entire east lot has been paved with either concrete or asphalt. The TAT could not verify that the east lot had been paved due to debris and piles of scrap metal covering the lot. A debris and ash pile was located in the area where the wire incinerator was formerly located. The sweat furnaces have been demolished. A foundation

of a building was located in the northeast corner of the east lot.

A small scale house is located on the west lot along the south side which is used to weigh incoming trucks loaded with scrap metal. A crane and 3 semi-trailers were parked along the west fence line of the west lot. Located along the north property line from west to east were a crane boom, a ash pile with tires placed on top, a striped out car, and 2 semi-trailers. Three large crates were located along the east fence line of the west lot.

During the reconnaissance CIEC workers were observed to burn a pile of cardboard material at the east lot and burn unknown materials in a 55-gallon drum at the west lot. Later CIEC workers put out the cardboard fire with water from a hose. The materials burned in the 55-gallon drum gave off a black smoke that was irritating to the eyes, nose, and throat. When the metal shredder was in operation a dust cloud could be observed. The metal shredder was turned off while TAT was in the area. Oil stains were observed on the ground at both lots. A motor had been cut open and oil was observed flowing from it at the east lot.

To characterize the possible hazardous substances on-site, the TAT collected ten soil samples including a duplicate sample and a matrix spike/matrix spike duplicate (MS/MSD) sample (see Figure 3 Sample Location Map). All soil samples were collected from 1 to 3 inches in depth with a shovel and a trowel.

All soil samples collected on-site were collected in Level D personal protection. Sampling gloves were changed before each sample was collected. Sampling equipment was decontaminated after each use, using a alconox and distilled water wash solution and triple rinsed with distilled water. All personal protective equipment was decontaminated as above and disposed of properly.

Soil samples S1, S2, S3, S7, and S10 were collected from the east lot. Soil sample S1 was collected from an area approximately 15 north of the northeast corner of the office. Soil Sample S2 was collected near a tree along the north property line west of the former wire incinerator. S3 was collected from a ash pile approximately 20 feet east of sample S2. An white crystalline substance was noted in the ash material and two dead rats were noted on top of the ash pile. Sample S7 was collected 25 feet north and 25 feet east of the northwest corner of the office building near the railroad tracks. Soil sample S10 was collected approximately 10 feet north and 35 feet east of the northwest corner of the office building.

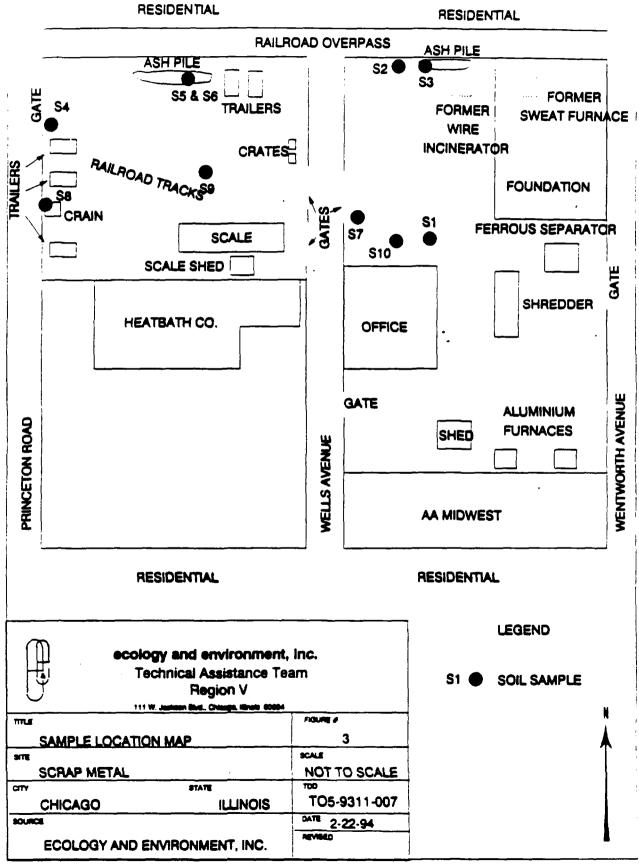


TABLE 3
ANALYTICAL DATA FOR PCBs
AT THE SCRAP METAL SITE
CHICAGO, ILLINOIS
unit = ppb

SAMPLE NUMBER	S1	S2	S3	S4	S5	S6	S 7	SB	S 9	S10
PARAMETER PCB 1016	77,000	ND	4,100	73,000	140,000	190,000	55,000	100,000	49,000	15,000
PCB 1254	45,000	ND_	18,000	170,000	510,000	300,000	42,000	550,000	83,000	24,000
PCB 1260	21,000	2,000,000	15,000	92,000	210,000	200,000	39,000	80,000	32,000	22,000
TOTAL PCBs	143,000	2,000,000	37,100	335,000	1,130,000	690,000	136,000	730,000	164,000	61,000

ND = Non Detect & your is Trisser.

1 Me Contraction of

TABLE 4 ANALYTICAL DATA FOR DIOXINS AT THE SCRAP METAL SITE CHICAGO, ILLINOIS

			UNCO, 121		·
Sample ID# Analyte	S1 µg/kg	52 µg/kg	S3 µg/kg	S8 μg/kg	BLANK µg/kg
TOTAL 2378-TCDD EQUIVALENCE	0.265	4.004	1.207	0.36	0
2378-TCDF	0.22	ND	1.7*	מא	ND
TOTAL TCDF	1.5	17	9.4	1.9	ND
2378-TCDD	ND	ND	ND	ND	סא
TOTAL TCDD	0.034	0.58	0.13	0.15	ND.
12378-PeCDF	0.094	4.4	0.55	ND	ND
23478-PeCDF	0.26	6.3	1.3	0.22	ND
TOTAL PeCDF	1.3	42	7.6	3.3	ND.
12378-PeCDD	מא	ND	ИD	ДИ	ИD
TOTAL PeCDD	ND	ND	ND	. ND	ND
123478-HxCDF	0.25	ND	0.81	0.5	סא
123678-HxCDF	0.16	ND	0.64	ND	ND
234678-HxCDF	0.23	2.7	0.99	0.36	0.039
123789-HxCDF	0.095	0.890	0.22	0.17	מא_
TOTAL HXCDF	2	15	7.7	2.7	0.039
123478-HxCDD	0.021*	ND	0.1	0.029	ИД
123678-HxCDD	0.070*	0.25	0.17	0.670	ND
123789-HxCDD	0.033*	0.13	0.12	0.19	מא
TOTAL HXCDD	0.24	0.38	1.1	1.8	ND
1234678-HCDF	0.9	12	3	1.1	ОИ
1234789-HpCDF	0.12	3.6	0.33	0.24	מא
TOTAL HPCDF	1.5	16	4.5	1.3	ND
1234678-HpCDD	0.73	2.9	1.5	2.8	0.0023
TOTAL HPCDD	1.4	5.9	2.9	5.7	0.0044
OCDF	0.77	30	1.6	1.6	ND
OCCD	4.3	22	3.9	51	0.021

ND - Non detect

^{* -} Value may include contributions from other TCDF isomers.

ppb. The PCBs can be directly associated with past activities at the site as reported by the Heatbath employee and former railroad employee. On-site soil contain dioxins that were greater than 1 ug/kg (2,3,7,8-TCDD equivalence) S2 4.004 ug/kg and S3 1.207 ug/kg. CIEC workers continue work on-site where they have the potential to come in contact with hazardous substances.

6.2 Migration of Contaminants

On-site soils contains heavy metals that were above the RCRA limits for cadmium and lead (1 and 5 mg/L, respectively) which are listed wastes D006 and D008, respectively. Total PCBs were detected in on-site soils ranging from 37,100 to 2,000,000 ppb. On-site soils contain dioxins that were greater than 1 ug/kg (2,3,7,8-TCDD equivalence) in samples S2 4.004 ug/kg and S3 1.207 ug/kg. The Agency for Toxic Substances and Disease Registry (ATSDR) considers 1 ug/kg of dioxin in soil to be a level of concern in residential areas. The migration of contaminants from the facility is possible due to piles by airborne dusts or storm run-off which could effect on-site workers and nearby residential areas.

7.0 SUMMARY

The site assessment documented the existence of hazardous substances at the Scrap Metal site. These substances were detected in samples of on-site soils which were found to contain materials that possess characteristics of hazardous wastes, as defined by RCRA. The SM site is an active metal reclamation facility. Conditions at the SM site may pose risks to the business employees on-site and to nearby residential areas.

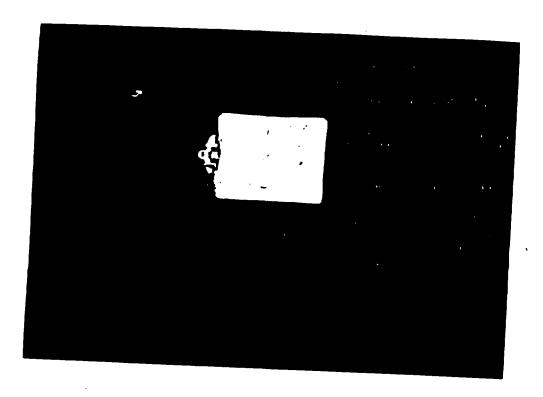
7.1 RCMS 4.0 COST PROJECTION

Removal cost estimates for off-site disposal were projected using the U.S. EPA Removal Cost Management System (RCMS) program. Cost estimates to mitigate threats due to on-site contaminated materials contained in the soil are based on the following assumptions:

- 1. There is approximately 4,500 cubic yards contaminated soils on-site. This estimated amount was generated assuming that the 3-acre site could be contaminated to a depth of at least 1 feet. This estimate does not include the soils underneath the concrete in east lot.
- Soil containing PCBs at concentrations grater than 50 ppm will be disposed of at a TSCA approved landfill;

- 3. Soil containing dioxin at concentrations grater than 1 ug/kg will be disposed of at a TSCA approved landfill;
- 4. Soil exhibiting TCLP cadmium and lead concentrations greater than 1 and 5 mg/l respectively will be stabilized either on-site or off-site and transported to an appropriate and RCRA permitted off-site landfill. If the soil also exhibits a PCB concentration greater than 50 mg/kg prior stabilization, the soil will be stabilized and disposed of at a TSCA approved landfill.

The cost estimates for the cleanup are based on 90 12-hour days by Emergency Response Cleanup Support (ERCS) contractor; Riedel Environmental Services, and is projected to cost \$1,696,050. See Appendix C for the RCMS cost estimate.



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1325
DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE
DESCR: SP COLLECTED FROM THE MIDDLE OF
THE WEST LOT BY RAILROAD TRACKS



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1250
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN MORDING

PHOTOGRAPHER: JOHN NORDINE DESCR: SB COLLECTED NEAR CRAIN ALONG



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1225
DIRECTION: HORTH

PHOTOGRAPHER: JOHN NORDINE
DESCR: S7 COLLECTED 25'N & 25'E OF THE
NW CORNER OF THE OFFICE



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1325 DIRECTION: NORTH
PHOTOGRAPHER: JOHN HORE



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1203
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT, S2 LOCATED ALONG
RAILROAD OVERPASS



SITE: SCRAP METAL
DATE: 2-22-94
TINE: 1205
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR- FAST LOT SOIL SAMPLE \$2

AND I



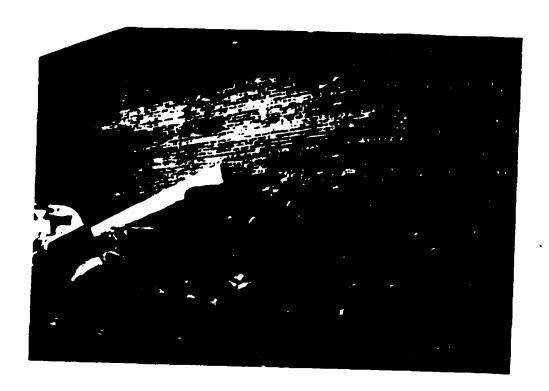
SITE: SCRAP METAL DATE: 2-22-94 1200 TIME:

DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE DESCR: \$1 EAST LOT, 20' NORTH OF THE OFFICES NE CORNER.



SITE: SCRAP METAL 2-22-94 DATE: TIME: DIRECTION: NORTH PHOTOGRAPHER: JOHN NORT W



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTH SIDE OF THE OFFICE
BUILDING EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1345 TIME: DIRECTION: SOUTH

100

Park Line

4. ----



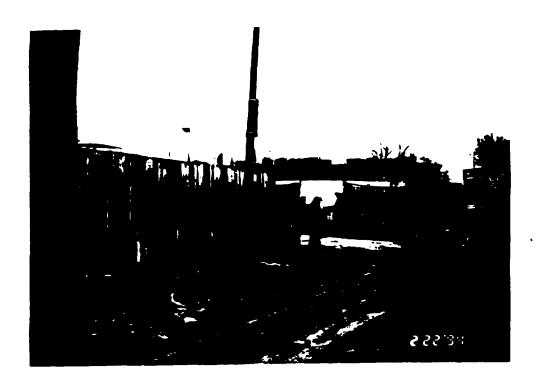
SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1225
DIRECTION: SOUTH

PHOTOGRAPHER: JOHN NORDINE

DESCR: OIL STAINED AREA WERE S10 WAS COLLECTED N SIDE OF THE OFFICE



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1225
DIRECTION: SOUTH
PHOTOGRAPHER- JOHN WORD! WE

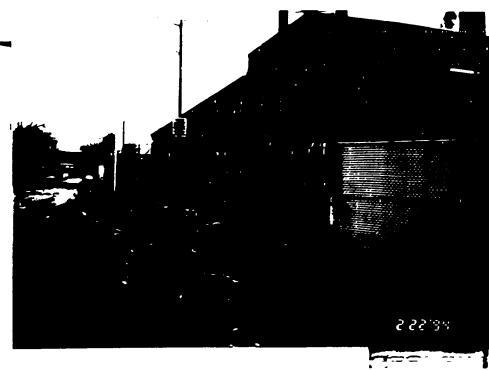


SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTH

PHOTOGRAPHER: JOHN NORDINE DESCR: WEST SIDE OF THE OFFICE NOTE FENCE



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1215 DIRECTION: NORTHWEST PHOTOGRAPHER: JOHN NORDINE DESCR. WORTHWEST AREA OF EAST 2"



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: NORTH
PHOTOGRAPHER: JOHN MORDINE
DESCR: FENCE ALONG THE WEST SIDE OF
THE EAST LOT

' .



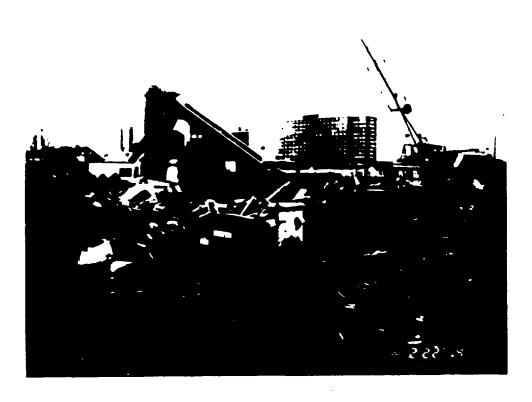
SITE: SCRAP METAL DATE: 2-22 2-22 % 1345 NOR!# TIME: DIRECTION:

PHOTOGRAPHER: JOHN WORD I ME



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDIN

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST SIDE OF THE OFFICE
BUILDING EAST LOT



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: MORTHEAST
DUOTOCCALDUED. HOW MORTHE

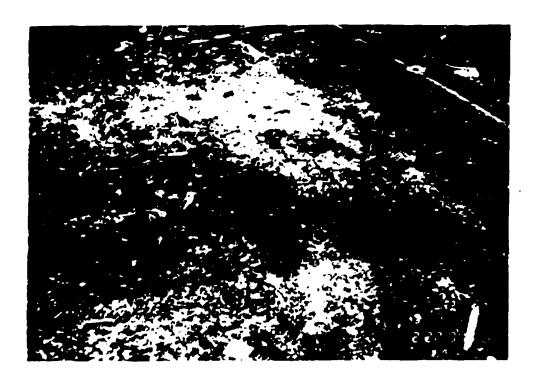


SITE: SCRAP METAL DATE: 2-22-94 TIME: 1230 TIME: DIRECTION: EAST

PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT CARDBOARD FIRE BY EAST
SIDE OF THE OFFICE BUILDING



SITE: SCRAP METAL 2-22-94 DATE: TIME: 1210 DIRECTION: \$0u*# PHOTOGRAPHER - JOHN MORO! WE



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1305
DIRECTION: NORTH
PHOTOGRAPHER: JOHN MORDINE

DESCR: S4 COLLECTED 3' EAST OF THE NORTHWEST GATE, WEST LOT

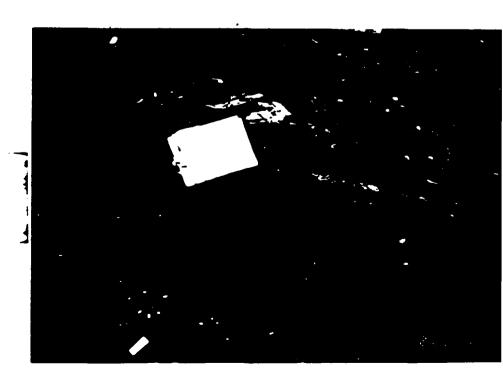


-SITE: SCRAP METAL
DATE: 2-22-96
TIME: 1310
DIRECTION: NORTH
PHOTOGRAPHER: JOHN MORDINE

SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1215
DIRECTION: NORTH
PHOTOCRAPHUR- JOHN NORDINE

到分。

DIRECTION: NORTH
PHOTOGRAPHER: JOHN NORDINE
DESCR: SOIL SAMPLE S3 FROM ASH PILE IN
THE EAST LOT, NOTE THE DEAD RATS





SITE: SCRAP HETAL 2-22-94 1345 DATE: TIME: DIRECTION:

PHOTOGRAPHER: JOHN NORDINE
DESCR: SCRAP METAL PILES EAST LOT NOTE
APARTMENTS IN THE BACKGROUND



SITE: SCRAP METAL 2-22-94 1205 DATE: TIME: DIRECTION: EAST

PHOTOGRAPHER: JOHN NORDINE

DESCR: EAST LOT SCRAP METAL PILES



SITE: SCRAP METAL DATE: 2-22-94
TIME: 1230
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN NORDINE

DESCR: EAST LOT RECIEVING SHED FOR RECYCLE MATERIALS



SITE: SCRAP DATE: TIME: 10411 DIRECTION:

PHOTOGRAPHER ACES MORDINE DESCR: DRUM WE ME CE IVING SHED



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTHEAST
PHOTOGRAPHER: JOHN HORDINE
DESCR: SOUTH PART OF THE EAST LOT



SITE: SCRAP METAL DATE: 2-22 TIME: 1345 2-22-94 1345

DIRECTION: EAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: RECYCLING SHED IN THE EAS' 2"



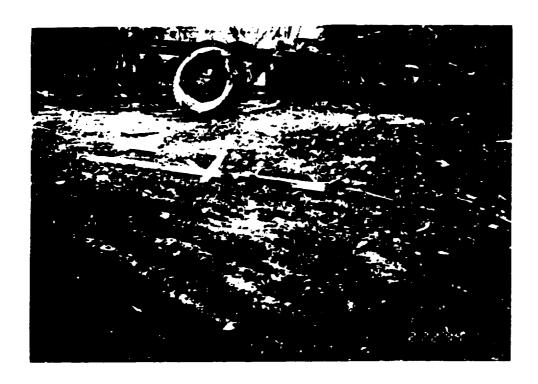
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1230 DIRECTION: WEST PHOTOGRAPHER: JOHN NORDINE

* **

DESCR: ALLMINUM FURNACES ALONG SOUTH PROPERTY LINE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94
TIME: 1345
DIRECTION: SOUTH
PHOTOGRAPHER: JOHN BORDINE
DESCR: SCRAP METAL PILES AND ALUMINUM



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1250
DIRECTION: NORTH

PHOTOGRAPHER: JOHN NORDINE
DESCR: CRAIN BOOM, ASH PILE, AND TIRES
ALONG NORTH SIDE OF WEST LOT



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1305
DIRECTION: EAST

DIRECTION: EAST PHOTOGRAPHER: JOHN NORDINE



SITE: SCRAP NETAL DATE: 2-22-94 TIME: 1250 DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: NORTHEAST PART OF THE WEST LOT



SITE: SCRAP METAL DATE: 2-22 TIME: 1250
DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN MCMD1ME



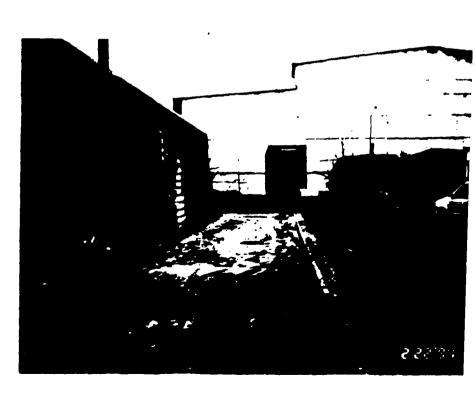
SITE: SCRAP METAL DATE: 2-22-94 TIME: 1300 DIRECTION: SOUTHWEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: SOUTHWEST PART OF THE WEST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1325 DIRECTION: NORTH



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1345
DIRECTION: NORTHWES DIRECTION: NORTHWEST
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST SIDE OF THE WEST LOT, NOTE
FENCE AND GATES



SITE: SCEME DATE: TIME: DIRECTION

PHOTOGRAPHO AND CORDINE HESTR. SCA. . MES MED TRUCK SCALE IN



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1225 DIRECTION: WEST
PHOTOGRAPHER: JOHN MORDINE
DESCR: METAL SHREDDER AT THE EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1330 DIRECTION: EAST



SITE: SCRAP METAL 2-22-94 1330 DATE: TIME: DIRECTION: NORTHEAST

PHOTOGRAPHER: JOHN NORDINE DESCR: EAST LOT SCRAP METAL PILES AND EMPTY DRUMS



SITE: SCRAP ME'AL DATE: 2 22 🛰 1365 TIME: DIRECTION: EAST

PHOTOGRAPHER: JOHN MORDINE
DESCR: FERROUS SOR'ER AND THE METAL
SHREDDER & 'ME EAST LOT



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1202 TIME: 1202 DIRECTION: NORTH PHOTOGRAPHER: JOHN NORDINE

DESCR: EAST LOT, PILES OF SCRAP METAL BY RAILROAD OVERPASS



SITE: SCRAP METAL DATE: 2-22-94 TIME: 1205 DIRECTION: NORTHEAST
PHOTOGRAPHER: JOHN NORDINE
DESCR: EAST LOT SCRAP METAL PILES '0



SITE: SCRAP METAL
DATE: 2-22-94
TIME: 1201
DIRECTION: MORTH
PHOTOGRAPHER: JOHN MORDINE

DESCR: EAST LOT RAILROAD OVERPASS NORTH BOUNDARY OF SITE.



SITE: SCRAP METAL
DATE: 2-22 %
TIME: 1202
DIRECTION: NORTH
PHOTOGRAPHER: JOHN #CMC #E

APPENDIX B DATA QUALITY ASSURANCE REVIEW



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111 Mest Jackson Boulevard On cago III noi**s** 60604 Teil 312 663-9415 (Faxil 312 663-019)

MENORANDU M

DATE: May 3, 1994

TO: John Nordine, Project Manager, E&E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, EGE, Chicago, IL NF

THRU: David Hendren, TAT-Chemist, E&E, Chicago, IL

SUBJ: Polychlorinated Biphenyl Data Quality Assurance Review, Scrap

Metal Site, Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 Project TDD: T059402007

Analytical PAN: EIL0831AAA Project PAN: EIL0831SAA

The data quality assurance review of 10 soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Polychlorinated Biphenyls (PCBs) was performed by Twin City Testing located in St. Paul. Minnesota, Illinois in accordance with U.S. EPA Method SW-846-8080.

The samples were numbered as following:

TAT Sample #	Corresponding to ->	Laboratory Sample #
S-1		15270
S-2		15277
s-3		15278
S-4		15279
S~5		15283
S ~6		15287
S-7		15289
S~8		15290
S-9		15291
S-10		15292

Data Qualifications:

I Sample Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/1/94, and analyzed on 3/6/94, and 3/7/94. The holding time criteria of 14 days for soil from collection to extraction was satisfied. The analysis of the samples was completed within the 40 day holding time requirement after extraction.

II Instrument Performance: Acceptable

The quality control criteria established for the surrogate retention time shift of less than 0.3% were acceptable. Peak resolution was adequate.

III Calibration: Qualified

A. Initial Calibration:

A 5-point calibration check was performed prior to sample analysis. The Relative Standard Deviations (RSDs) of all calibration factors for all Arochlors did not meet the quality control criteria of less than or equal to 10% RSD except 1260 on the DB-608 column, therefore; positive results were flagged (J) as estimated.

B. Continuing Calibration:

The percent difference (%D) requirements of less than 15% were acceptable for all data.

IV Method Blanks: Acceptable

A method blank was analyzed with the samples. No contaminants were detected above the instrument detection limit.

V Matrix Spike/Matrix Spike Duplicate: Acceptable

Spike Sample Analysis:

All Matrix Spike/Matrix Spike Duplicate recoveries were diluted out due to high level of PCBs in the samples. No action was necessary.

VI Compound Identification and Quantitation: Acceptable

All reported Retention Times were within the allowed RT windows. Positive results were confirmed using dual column analysis. The laboratory states in the narrative that because of the complexity of the Aroclor patterns present, it is possible that several Aroclors (PCB 1232 and PCB 1242) could be misidentified as PCB 1016.

VII Surrogate Recovery: Acceptable

The surrogates were diluted out due to high level of PCBs in the samples. No action was necessary.

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality control Guidance For Removal Activities" (OSWER Directive 9360.4-01, April 1990).

Based upon the information provided, the data are acceptable for use with the above stated data qualifications.

Data Qualifiers and Definitions

J - The associated numerical value is an estimated quantity because the were less than the contract required detection limits or quality control criteria were not met.

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-1

TCT ID:

15270

ICT ID:	152/0		
Parameter:		POL	
PCB 1016	77,000	7,800	
PCB 1221	ND ·	7,800	
PCB 1232	ND	7,800	
PCB 1242	ND	7,800	
PCB 1248	ND	7,800	
PCB 1254	45,000	7,800	
PCB 1260	21,000	7,800	
% Surrogate #1 Recovery:	1%		
% Surrogate #2 Recovery:	—¹ %	•	

Date Extracted:

3/1/94

Date Analyzed:

3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition.

¹Surrogate diluted out due to high level of PCB in sample.

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-2

TCT ID:

15277

Parameter:		POL
PCB 1016	ND	250,000
PCB 1221	ND .	250,000
PCB 1232	ND	250,000
PCB 1242	ND .	250,000
PCB 1248	ND	250,000
PCB 1254	ND	250,000
PCB 1260	2,000,000	250,000
% Surrogate #1 Recovery:	¹ %	

% Surrogate #2 Recovery:

--¹ %

Date Extracted:

3/1/94

Date Analyzed:

3/7/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in ug/Kg which is equal to parts-per-billion)

Client ID:

S-3

TCT ID:	15278	
Parameter:		POL
PCB 1016	4,100	2,500
PCB 1221	ND ·	2,500
PCB 1232	ND	2,500
PCB 1242	ND	2,500
PCB 1248	ND	2,500
PCB 1254	18,000	2,500
PCB 1260	15,000	2,500
% Surrogate #1 Recovery:	-1%	
% Surrogate #2 Recovery:	-1%	
Date Extracted:	3/1/94	
Date Analyzed:	3/7/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

POL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachioro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference: EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-4

TCT ID: 15279

Parameter:		POL
PCB 1016	73,000	64,000
PCB 1221	ND ·	64,000
PCB 1232	ND	64,000
PCB 1242	ND	64,000
PCB 1248	ND	64,000
PCB 1254	160,000	. \$64,000
PCB 1260	92,000	64,000
% Surrogate #1 Recovery:	-¹ %	
% Surrogate #2 Recovery:	-1 %	

Date Extracted:

3/1/94

Date Analyzed:

3/6/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-5

TCT ID:	15283	
Parameter:		POL
PCB 1016	410,000	75,000
PCB 1221	ND .	75,000
PCB 1232	ND	75,000
PCB 1242	ND	75,000
PCB 1248	ND	75,000
PCB 1254	510,000	75,000
PCB 1260	210,000	75,000
% Surrogate #1 Recovery:	¹ %	
% Surrogate #2 Recovery:	¹%	
		•
Date Extracted:	3/1/94	

3/6/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-6

15287 TCT ID:

Parameter:		POL
PCB 1016	190,000	72,000
PCB 1221	ND	72,000
PCB 1232	ND	72,000
PCB 1242	ND .	72,000
PCB 1248	ND	72,000
PCB 1254	300,000	72,000
PCB 1260	200,000	72,000
% Surrogate #1 Recovery:	_1%	
% Surrogate #2 Recovery:	¹ %	

Date Extracted:

3/1/94

Date Analyzed:

3/6/94

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachioro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition. Reference:

¹Surrogate diluted out due to high level of PCB in sample.

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-7

TCT ID:	15289	
Parameter:		POL
PCB 1016	55,000	13,000
PCB 1221	ND	13,000
PCB 1232	ND	13,000
PCB 1242	ND	13,000
PCB 1248	ND	13,000
PCB 1254	42,000	13,000
PCB 1260	39,000	13,000
% Surrogate #1 Recovery:	1 %	
% Surrogate #2 Recovery:	¹ %	
Date Extracted:	3/1/94	•

3/7/94

All results are reported on a dry weight basis.

POL = Practical Quantitation Limit

ND = Not Detected

Date Analyzed:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



¹Surrogate diluted out due to high level of PCB in sample.

(All values are in µg/Kg which is equal to parts-per-billion)

	lient	m.
L. 1	исии	ID:

S-8

TCT ID:	15290		
Parameter:		POL	
PCB 1016	100,000²	75,000	
PCB 1221	ND	75,000	
PCB 1232	ND	75,000	
PCB 1242	ND	75,000	
PCB 1248	ND	75,000	
PCB 1254	550,000	75,000	
PCB 1260	80,000	75,000	
% Surrogate #1 Recovery:	¹ %		
% Surrogate #2 Recovery:	¹ %		
Date Extracted:	3/1/94		
Date Analyzed:	3/7/94		

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition.



²Value was not confirmed within 25%, the lower value is reported.

(All values are in µg/Kg which is equal to parts-per-billion)

S-9

Client ID:

TCT ID:	15291	
Parameter:		POL
PCB 1016	49,000	14,000
PCB 1221	ND ·	14,000
PCB 1232	ND	14,000
PCB 1242	ND	14,000
PCB 1248	ND	14,000
PCB 1254	83,000	14,000
PCB 1260	32,000	14,000
% Surrogate #1 Recovery:	1 %	
% Surrogate #2 Recovery:	¹ %	
Date Extracted:	3/1/94	
Date Analyzed:	3/7/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

Huntingdon

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-10

TCT ID:	15292	
Parameter:		POL
PCB 1016	15,000	5,000
PCB 1221	ND ·	5,000
PCB 1232	ND	5,000
PCB 1242	ND	5,000
PCB 1248	ND	5,000
PCB 1254	24,000	5,000
PCB 1260	22,000	5,000
% Surrogate #1 Recovery:	-1%	
% Surrogate #2 Recovery:	_¹ %	
		•
Date Extracted:	3/1/94	
Date Analyzed:	3/7/94	

¹Surrogate diluted out due to high level of PCB in sample.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate #1 = TCMX (2,4,5,6-tetrachloro-m-xylene)

Surrogate #2 = DCB (decachlorobiphenyl)

Reference:

EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition.





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111 Mest Jackson Boulevard On pago I Inois 60604 Tel: 310 663-9415 Fax: 312 663-0791

MEMORANDUM

DATE: May 3, 1994

TO: John Nordine, Project Manager, E&E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, E&E, Chicago, IL Nf

THRU: David Hendren, TAT-Chemist, E&E, Chicago, IL

SUBJ: Inorganic Data Quality Assurance Review, Scrap Metal Site,

Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 Project TDD: T059402007 Analytical PAN: EIL0831AAA Project PAN: EIL0831SAA

The data quality assurance review of 10 Soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Total and TCLP Metals was performed by Twin City Testing located in

St. Paul, Minnesota in accordance with U.S. EPA Methods SW-846-6010/7000.

The samples were numbered as following:

TAT Sample #	Corresponding to ->	Laboratory Sample #.
S-1		15270
S-2		15277
s-3		15278
s-4		15279
s-5		15283
s -6		15287
S-7		15289
S -8		15290
s -9		15291
S-10		15292

Data Qualifications:

I Sample Holding Time: Acceptable

The samples were collected on 2/22/94 and analyzed between 3/1/94, and 3/18/94. The holding time criteria of 6 months for metals and 28 days for mercury from collection to analysis was satisfied.

II Calibration: Acceptable

A. Initial Calibration:

Calibration results were within the established quality control limits of 90-110% of the true value for metals, and 80-120% for mercury.

B. Continuing Calibration:

Calibration results showed that the control criteria of 90-110% for metals, and 80-120% for mercury were satisfied.

III Method Blanks: Acceptable

A method blank was analyzed with the samples. No contaminants were detected above the instrument detection limit.

IV Interference Check Sample Analysis: Acceptable

All parameters were within the Interference Check Sample (ICS) control limits of 80-120% of the true values. ICS was run at the beginning and end of sample analysis.

V Matrix Spike/Matrix Spike Duplicate:

Spike Sample Analysis: Acceptable

All Matrix Spike/Matrix Spike Duplicate recoveries were within the control limits of 80-120%.

VI Optional Additional QC:

Laboratory Control Sample Analysis: Acceptable

The quality control criteria of 80-120% were met for the control sample.

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance For Removal Activities" (OSWER Directive 9360.4-01, April 1990).

Based upon the information provided, the data are acceptable for use.

(All values are in µ2:L which is equivalent to parts-per-billion)

Client ID:

SI

S2

S3

TCT ID:

15270

15277

15278

<u>Parameter</u>				POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	ND	100	3/15/94	6010
Barium	1,500	2,800	2,700	10	3/15/94	6010
Cadmium	870	1,300	110	10	3/15/94	6010
Chromium	ND	ND	ND	10	3/15/94	6010
Lead	7,500	16,000	690	50	3/15/94	6010
Mercury	ND	ND	ND	0.40	3/18/94	7470
Selenium	ND	ND	ND	100	3/15/94	6010
Silver	ND	ND	ND	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94	·		

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste. SW-846, November 1986, 3rd Edition.

(All values are in µg/L which is equivalent to parts-per-billion, unless noted otherwise)

Client ID:

S4

S4

S4

15279D 15279S 15279 TCT ID: Test Test **POL Parameter Date Method** 100 3/15/94 ND 106% 6010 Arsenic ND 80% 10 3/15/94 6010 1,100 1,100 Barium 220 90% 10 3/15/94 6010 Cadmium 230 10 3/15/94 6010 Chromium 10 ND 91% 50 3/15/94 6010 8,100 81% Lead 8,200 125% 0.40 3/18/94 7470 ND ND Mercury 100 3/15/94 6010 ND 113% Selenium ND ND 92% 10 3/15/94 6010 Silver ND TCLP Date: 3/14/94 3/14/94 3/14/94

PQL = Practical Quantitation Limit

ND = Not Detected

D = Duplicate

S = Spike

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

(All values are in $\mu g/L$ which is equivalent to parts-per-billion)

Client ID:

S5

S6

S7

TCT ID:

15283

15287

15289

<u>Parameter</u>				POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	ND	100	3/15/94	6010
Barium	4,800	2,700	1,300	10	3/15/94	6010
Cadmium	910	660	1,200	10	3/15/94	6010
Chromium	ND	ND	ND	10	3/15/94	6010
Lead	45,000	4,300	7,400	50	3/15/94	6010
Mercury	ND	ND	ND	0.40	3/18/94	7470
Selenium	ND	ND	ND	100	3/15/94	6010
Silver	ND	ND	ND	10	3/15/94	6010
TCLP Date:	3/14/94	3/14/94	3/14/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

LABORATORY NO: 4416 94-2757

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(All values are in µg/L which is equivalent to parts-per-billion)

Client ID:

S8

S9

S10

TCT ID:

15290

15291

15292

Parameter				POL	Test <u>Date</u>	Test <u>Method</u>	
Arsenic	ND	ND	ND	100	3/15/94	6010	
Barium	1,500	1,600	2,000	10	3/15/94	6010	
Cadmium	850	480	690	10	3/15/94	6010	
Chromium	12	15	ND	10	3/15/94	6010	
Lead	5,300	71,000	32,000	50	3/15/94	6010	
Mercury	ND	ND	ND	0.40	3/18/94	7470	
Selenium	ND	ND	ND	100	3/15/94	6010	
Silver	ND	ND	ND	, 10	3/15/94	6010	
TCLP Date:	3/14/94	3/14/94	3/14/94				

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.



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MEMORANDUM

DATE: March 29, 1994

TO: John Nordine, Project Manager, E & E, Chicago, IL

FROM: Nabil Fayoumi, TAT-Chemist, E & E, Chicago, IL

THRU: David Hendren, TAT-Chemist, E & E, Chicago, IL

SUBJ: Organic Data Quality Assurance Review, Scrap Metal Site,

Chicago, Cook County, Illinois.

REF: Analytical TDD: T059402807 Project_TDD: T059402007

Analytical PAN: EILO831AAA Project PAN: EILO831SAA

The data quality assurance review of 10 soil samples collected from the Scrap Metal Site in Chicago, Illinois has been completed. Analysis for Semivolatile Organics (SVOA) was performed by Twin City Testing located in St. Paul, Minnesota in accordance with U.S. EPA SW-846 method 8270.

The samples were numbered as following:

TAT Sample #	Corresponding to ->	Laboratory Sample #
S-1		15270
S-2		15277
s-3		15278
S-4		15279
S-5		15283
s-6		15287
5-7		15289
5-8		15290
5-9		15291
S-10		15292

Data Qualifications:

I Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/3/94, and analyzed between 3/8/94 and 3/10/94. The holding time criteria from collection to extraction of 14 days was satisfied. Extracts were analyzed within 40 days.

II GC/MS Tuning: Acceptable

GC/MS ion abundance criteria using Decafluorotriphenylphosphine (DFTPP) were acceptable.

III Calibration: Acceptable

A. Initial Calibration:

A 5-point initial calibration was performed prior to analysis. All average relative response factors were greater than 0.05. All percent relative standard deviation (%RSD) between response factors were less than 30%, except 2,4-dinitrophenol (33.1 % RSD), which was not detected in any sample; therefore qualification was not necessary.

B. Continuing Calibration:

The percent difference (%D) between initial and continuing calibration were within the quality control criteria of less than or equal to 25% for all detected compounds.

IV Method Blank: Acceptable

The method blank was contaminated with Di-n-butylphthalate (47 PPS). No action was necessary because all reported sample values were greater than 10 times blank concentration level.

V Surrogate Recovery: Acceptable

Surrogates in the matrix spike sample (number S-4MS) were recovered at less than 10% recovery. Since both samples S-4 and S-4MSD displayed acceptable surrogate recoveries, it appears that the poor surrogate response in sample S-4MS resulted from sample preparation problems.

VI Matrix Spike/Matrix Spike Duplicates: Not Applicable

Fourteen out of 22 percent recoveries and 10 out of 11 relative percent differences for the MS/MSD were outside the established quality control criteria. This again probably resulted from sample S-4MS preparation problems.

VII Internal Standards: Qualified

The internal standard (chrysene D-12) area counts were outside (below) the range of -50% to +100% from the associated calibration standard for the following samples S-3, S-4, S-5, S-6, S-7, S-8, S-9, and S-10; therefore, all positive values quantitated from chrysene D-12 in these samples are flagged "J" (estimated).

VIII Overall Assessment of Data for Use

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance for Removal Activities" (OSWER 9360.4-01 April, 1990). Sased upon the information provided, the data are acceptable for use with the above stated qualifications.

Data Qualifiers and Definitions:

J - The associated numerical value is an estimated quantity because the reported concentration is less than the contract required detection limit or quality control criteria were not met.

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S1 Lab ID (HSN): 15270
Matrix: SOIL Filename: 4068K08

Date Sampled: 02/22/94 Sample Size: 15.5 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 24.08

Compounds:	ug/Kg (PPB)	EQL
Phenol	1700	1700
bis(2-Chloroethyl)ether	1700 U	1700
2-Chlorophenol	1700 U	1700
1,3-Dichlorobenzene	1700 U	1700
1,4-Dichlorobenzene	1700 U	1700
1,2-Dichlorobenzene	1700 U	1700
2-Methylphenol	1700 U	1700
2,2'-oxybis(1-Chloropropane)	1700 U	1700
4-Methylphenol	220 J	1700
N-Nitroso-di-n-propylamine	1700 U	1700
Hexachloroethane	1700 U	1700
Nitrobenzene	1700 U	1700
Isophorone	1700 U	1700
2-Nitrophenol	1700 U	1700
2,4-Dimethylphenol	1700 U	1700
bis(2-Chloroethoxy)methane	1700 U	1700
2,4-Dichlorophenol	1700 U	1700
1,2,4-Trichlorobenzene	1300 J	1700
Naphthalene	970 J	1700
4-Chloroaniline	' 1700 Ŭ	1700
Hexachlorobutadiene	1700 U	1700
4-Chloro-3-methylphenol	1700 U	1700
2-Methylnaphthalene	2900	1700
Hexachlorocyclopentadiene	1700 U	1700
2,4,6-Trichlorophenol	1700 U	1700
2,4,5-Trichlorophenol,	4200 U	4200
2-Chloronaphthalene	1700 U	1700
2-Nitroaniline	4200 U	4200
Dimethylphthalate	510 J	1700
Acenaphthylene	1700 U	1700
2,6-Dinitrotoluene	1700 U	1700
3-Nitroaniline	4200 U	4200
Acenaphthene	710 J	1700
2,4-Dinitrophenol	4200 U	4200
4-Nitrophenol	4200 U	4200
Dibenzofuran	1700 U	1700
2,4-Dinitrotoluene	1700 U	1700
Diethylphthalate	1700 U	1700
4-Chlorophenyl-phenylether	1700 U	1700
Fluorene	1100 J	1700
4-Nitroaniline	4200 U	4200
4,6-Dinitro-2-methylphenol	4200 U	4200
(continued)		

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S1
Matrix: SOIL

Lab ID (HSN): 15270 Filename: 4068K08

	 	
Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine 4-Bromophenyl-phenylether Hexachlorobenzene Pentachlorophenol Phenanthrene Anthracene	1700 U 1700 U 1700 U 4200 U 3900 290 J	1700 1700 1700 4200 1700
Carbazole Di-n-butylphthalate Fluoranthene Pyrene Butylbenzylphthalate	1700 U 2600 B 2500 4200 2500	1700 1700 1700 1700 1700
3,3'-Dichlorobenzidine Benz(a)anthracene Chrysene	1700 U 1700 J 1900	1700 1700 1700
<pre>bis(2-Ethylhexyl)phthalate Di-n-octylphthalate Benzo(b)fluoranthene Benzo(k)fluoranthene Benzo(a)pyrene</pre>	100000 E 1700 U 2800 1100 J - 2000	1700 1700 1700 1700 1700
Indeno(1,2,3-cd) pyrene Dibenz(a,h) anthracene Benzo(g,h,i) perylene	1600 J 1700 U 1900	1700 1700 1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	55 %	25-121%
Phenol-d5	63 %	24-113*
2-Chlorophenol-d4	76%	20-130%
1,2-Dichlorobenzene-d4	87%	20-130%
Nitrobenzene-d5	51%	23-120%
2-Fluorobiphenyl	114%	30-115%
2,4,6-Tribromophenol	57%	19-122
Terphenyl-d14	112*	18-137*

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Compounds:	ug/Kg (PPB)	EQL
Phenol	3100 JD	17000
bis(2-Chloroethyl)ether	17000 UD	17000
2-Chlorophenol	17000 UD	17000
1,3-Dichlorobenzene	17000 UD	17000
1,4-Dichlorobenzene	17000 UD	17000
1,2-Dichlorobenzene	17000 UD	17000
2-Methylphenol	17000 UD	17000
2,2'-oxybis(1-Chloropropane)	17000 UD	17000
4-Methylphenol	17000 UD	17000
N-Nitroso-di-n-propylamine	17000 UD	17000
Hexachloroethane	17000 UD	17000
Nitrobenzene	17000 UD	17000
Isophorone	.17000 UD	17000
2-Nitrophenol	17000 UD	17000
2,4-Dimethylphenol	17000 UD	17000
bis(2-Chloroethoxy)methane	17000 UD	17000
2,4-Dichlorophenol	17000 UD	17000
1,2,4-Trichlorobenzene	17000 UD	17000
Naphthalene	17000 UD	17000
4-Chloroaniline	17000 UD	17000
Hexachlorobutadiene	17000 UD	17000
4-Chloro-3-methylphenol	17000 UD	17000
2-Methylnaphthalene	3200 JD	17000
Hexachlorocyclopentadiene	17000 UD	17000
2,4,6-Trichlorophenol	17000 UD	17000
2,4,5-Trichlorophenol	42000 UD	42000
2-Chloronaphthalene	17000 UD	17000
2-Nitroaniline	42000 UD	42000
Dimethylphthalate	17000 UD	17000
Acenaphthyl ene	17000 UD	17000
2,6-Dinitrotoluene	17000 UD	17000
3-Nitroaniline	42000 UD	42000
Acenaphth ene	17000 UD	17000
2,4-Dinitrophenol	42000 UD	42000
4-Nitrophenol	42000 UD	42000
Dibenzofuran	17000 UD	17000
2,4-Dinitrotoluene	17000 UD	17000
Diethylphthalate	17000 UD	17000
4-Chlorophenyl-phenylether	17000 UD	17000
Fluorene	1700 JD	17000
4-Nitroaniline	42000 UD	42000
4,6-Dinitro-2-methylphenol	42000 UD	42000

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS CONTINUED: EPA METHOD 8270

Client ID: S1
Matrix: SOIL

Lab ID (HSN): 15270 DL Filename: 4067L15

Compounds:	ug/Kg (PI	PB)	EQL
N-Nitrosodiphenylamine	17000	נים כים	17000
4-Bromophenyl-phenylether	17000	ים כים	17000
Hexachlorobenzene	17000		17000
Pentachlorophenol	42000	UD 4	42000
Phenanthrene	4100	JD :	17000
Anthracene	17000		17000
Carbazole	17000		17000
Di-n-butylphthalate	4700		17000
Fluoranthene	2900		17000
Pyrene	7100		17000
Butylbenzylphthalate	7700	JD 3	17000
3,3'-Dichlorobenzidine	17000	UD 1	17000
Benz(a)anthracene	1800	JD 1	17000
Chrysene	2000	JD 1	17000
bis(2-Ethylhexyl)phthalate	200000	ED 1	17000
Di-n-octylphthalate	17000	UD 1	17000
Benzo(b) fluoranthene	3000	JD 1	17000
Benzo(k) fluoranthene	17000	UD 1	17000
Benzo(a)pyrene	. 1900	JD 1	17000
Indeno(1,2,3-cd)pyrene	17000	מס מי	17000
Dibenz(a,h)anthracene	17000	ני סט	17000
Benzo(g,h,i)perylene	17000	ני סט	17000

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl 2,4,6-Tribromophenol	79%JD 93%JD 89%JD 90%JD 90%JD 116%JD 73%JD	QC LIMITS 25-121% 24-113% 20-130% 20-130% 23-120% 30-115% 19-122%
2,4,6-Tribromophenol Terphenyl-d14		-

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a try weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edit: :

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S1 Lab ID (HSN): 15270 DL2
Matrix: SOIL Filename: 4068K04

Date Sampled: 02/22/94 Sample Size: 15.5 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 20
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 24.08

Compounds:	ug/Kg	(PPB)	EQL
Phenol	340	ാ സ	34000
bis(2-Chloroethyl)ether	340	00 UD	34000
2-Chlorophenol	340	OO UD	34000
1,3-Dichlorobenzene	340	00 UD	34000
1,4-Dichlorobenzene	340	00 UD	34000
1,2-Dichlorobenzene	340	00 UD	34000
2-Methylphenol	340	00 UD	34000
2,2'-oxybis(1-Chloropropane)	340	00 UD	34000
4-Methylphenol	340	00 UD	34000
N-Nitroso-di-n-propylamine	340	00 UD	34000
Hexachloroethane	340	00 W	34000
Nitrobenzene	340	00 UD	34000
Isophorone	340	00 UD	34000
2-Nitrophenol	-340	00 UD	34000
2,4-Dimethylphenol	340	00 UD	34000
bis(2-Chloroethoxy)methane	340	00 UD	34000
2,4-Dichlorophenol	340	00 UD	34000
1,2,4-Trichlorobenzene	340	00 UD	34000
Naphthalene	340	00 UD	34000
4-Chloroaniline	340	00 W	34000
Hexachlorobutadiene	340	00 UD	34000
4-Chloro-3-methylphenol	340	00 UD	34000
2-Methylnaphthalene	340	00 UD	· 34000
Hexachlorocyclopentadiene	340	QU 00	34000
2,4,6-Trichlorophenol	340	00 UD	34000
2,4,5-Trichlorophenol		00 UD	
2-Chloronaphthalene		00 D	
2-Nitroaniline		00 D	
Dimethylphthalate		00 DD	
Acenaphthylene		00 UD	
2,6-Dinitrotoluene		00 Œ	
3-Nitroaniline		0,0 AD	
Acenaphth ene		00 m	
2,4-Dinitrophenol		00 D	
4-Nitrophenol		00 W	85000
Dibenzofuran		00 W	
2,4-Dinitrotoluene		00 MD	
Diethylphthalate		00 UD	
4-Chlorophenyl-phenylether		00 UD	
Fluorene		00 m	
4-Nitroaniline		00 UD	
4,6-Dinitro-2-methylphenol	850	00 D	85000 -

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S1
Matrix: SOIL

Lab ID (HSN): 15270 DL2 Filename: 4068K04

Compounds:	ug/Kg (P	PB)	EQL
N-Nitrosodiphenylamine	34000	סט	34000
4-Bromophenyl-phenylether	34000	UD	34000
Hexachlorobenzene	34000		34000
Pentachlorophenol	85000		85000
Phenanthrene	4600		34000
Anthracene	34000		34000
Carbazole	34000		34000
Di-n-butylphthalate	4900		34000
Fluoranthene	3700		34000
Pyrene		Ĵ	
Butylbenzylphthalate		JD	
3,3'-Dichlorobenzidine	34000		
Benz(a)anthracene	34000		
Chrysene	34000		
bis(2-Ethylhexyl)phthalate	180000		34000
Di-n-octylphthalate	34000	UD	34000
Benzo(b)fluoranthène	34000		34000
Benzo(k)fluoranthene	34000	\mathbf{w}	34000
Benzo(a) pyrene	34000		34000
Indeno(1,2,3-cd)pyrene	34000	ϖ	34000
Dibenz(a,h)anthracene	34000		34000
Benzo(g,h,i)perylene	34000	UD	34000

Surrogate Recovery 2-Fluorophenol	0 E \$ TD	QC LIMITS
•	85 %J D	25-121%
Phenol-d5	101 % JD	24-113%
2-Chlorophenol-d4	91 % JD	20-130%
1,2-Dichlorobenzene-d4	92 % JD	20-130%
Nitrobenzene-d5	108 %J D	23-120%
2-Fluorobiphenyl	112 % JD	30-115%
2,4,6-Tribromophenol	75 %J D	19-122%
Terphenyl-d14	168 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Eva. Lating Solid Waste", SW-846,

November 1986, 3rd Edit::n

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S2 Lab ID (HSN): 15277 Matrix: SOIL Filename: 4068K09

Date Sampled: 02/22/94
Date Received: 02/24/94
Date Extracted: 03/03/94
Date Analyzed: 03/09/94 Sample Size: 15.2 grams
Extract Vol.: 1000 uL
Dil. Factor: 1
GPC Factor: 2 GPC Factor: 2

% Moisture: 28.24

Compounds:	ug/Kg (PPB)	EQL
Phenol	1800 U	1800
bis(2-Chloroethyl)ether	1800 U	1800
2-Chlorophenol	1800 U	1800
1,3-Dichlorobenzene	1800 U	1800
1,4-Dichlorobenzene	1800 U	1800
1,2-Dichlorobenzene	1800 U	´1800
2-Methylphenol	1800 U	1800
2,2'-oxybis(1-Chloropropane)	1800 U	1800
4-Methylphenol	1800 U	1800
N-Nitroso-di-n-propylamine	1800 U	1800
Hexachloroethane	1800 U	1800
Nitrobenzene	1800 U	1800
Isophorone	1800 U	1800
2-Nitrophenol	- 1800 U	1800
2,4-Dimethylphenol	1800 U	1800
bis(2-Chloroethoxy)methane	1800 U	1800
2,4-Dichlorophenol	1800 U	1800
1,2,4-Trichlorobenzene	7500	1800
Naphthalene	180 J	1800
4-Chloroaniline	1800 U	1800
Hexachlorobutadiene	1800 U	1800
4-Chloro-3-methylphenol	1800 U	1800
2-Methylnaphthalene	180 J .	
Hexachlorocyclopentadiene	1800 U	1800
2,4,6-Trichlorophenol	1800 U	1800
2,4,5-Trichlorophenol	4600 U	4600
2-Chloronaphthalene	1800 U	1800
2-Nitroaniline	4600 U	4600
Dimethylphthalate	1800 U	1800
Acenaphthylene	1800 U	1800
2,6-Dinitrotoluene	1800 U	1800
3-Nitroaniline	4600 U	4600
Acenaphthene	1800 U	1800
2,4-Dinitrophenol	4600 U	4600
4-Nitrophenol	4600 U	4600 1800
Dibenzofuran	1800 U 1800 U	1800
2,4-Dinitrotoluene		1800
Diethylphthalate	1800 U 1800 U	1800
4-Chlorophenyl-phenylether	1800 U	1800
Fluorene	4600 U	4600
4-Nitroaniline	4600 U	4600
4,6-Dinitro-2-methylphenol	4000 0	4000

(continued)

TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS (CONTINUED) EPA METHOD 8270

Client ID: S2
Matrix: SOIL

Lab ID (HSN): 15277 Filename: 4068K09

ug/Kg (PPB)	EQL
1800 U	1800
1800 U	1800
1800 U	1800
4600 U	4600
1300 J	1800
180 J	1800
1800 U	1800
2200 B	1800
1200 J	1800
2100	1800
1800 U	1800
1800 U	1800
1100 J	. 1800
	1800
	1800
	1800
_	1800
	1800
	1800
	1800
	1800
1700 J	1800
	1800 U 1800 U 1800 U 4600 U 1300 J 180 J 1800 U 2200 B 1200 J 2100 1800 U

Surrogate Recovery		QC LIMITS
2-Fluorophenol	36 %	25-121%
Phenol-d5	38%	24-113%
2-Chlorophenol-d4	47%	20-130%
1,2-Dichlorobenzene-d4	56 %	20-130%
Nitrobenzene-d5	33 % J	23-120%
2-Fluorobiphenyl	61%	30-115%
2,4,6-Tribromophenol	32%	19-122%
Terphenyl-d14	70%	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

= Also detected in the associated Blank

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

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TCL SEMIVOLATILE ORGANIC COMPOUND RESULTS EPA METHOD 8270

Client ID: S2 Lab ID (HSN): 15277 DL Matrix: SOIL Filename: 4067L16

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Received: 02/24/94 Extract Vol.: 1000
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 28.24

Compounds:	ug/Kg (PPB)	EQL
Phenol	18000 UD	18000
bis(2-Chloroethyl)ether	18000 UD	18000
2-Chlorophenol	18000 UD	18000
1,3-Dichlorobenzene	18000 UD	18000
1,4-Dichlorobenzene	18000 UD	18000
1,2-Dichlorobenzene	18000 UD	18000
2-Methylphenol	18000 UD	18000
2,2'-oxybis(1-Chloropropane)	18000 UD	18000
4-Methylphenol	18000 UD	18000
N-Nitroso-di-n-propylamine	18000 UD	18000
Hexachloroethane	18000 UD	18000
Nitrobenzene	18000 UD	18000
Isophorone	18000 UD	18000
2-Nitrophenol	18000 UD	18000
2,4-Dimethylphenol	18000 UD	18000
bis(2-Chloroethoxy)methane	18000 UD	18000
2,4-Dichlorophenol	18000 UD	18000
1,2,4-Trichlorobenzene	6000 JD	18000
Naphthalene	18000 UD	18000
4-Chloroaniline	18000 UD	18000
Hexachlorobutadiene	18000 UD	18000
4-Chloro-3-methylphenol	18000 UD	18000
2-Methylnaphthalene	18000 UD	18000
Hexachlorocyclopentadiene	18000 UD	18000
2,4,6-Trichlorophenol	18000 UD	18000
2,4,5-Trichlorophenol	46000 UD	46000
2-Chloronaphthalene	18000 UD	18000
2-Nitroaniline	46000 UD	46000
Dimethylphthalate	18000 UD	18000
Acenaphthyl ene	18000 UD	18000
2,6-Dinitrotoluene	18000 UD	18000
3-Nitroaniline	46000 UD	46000
Acenaphthene	18000 UD	18000
2,4-Dinitrophenol	46000 UD	46000
4-Nitrophenol	46000 UD	46000
Dibenzofuran	18000 UD	18000
2,4-Dinitrotoluene	18000 UD	18000
Diethylphthalate	18000 UD	18000
4-Chlorophenyl-phenylether	18000 UD	18000
Fluorene	18000 UD	18000
4-Nitroaniline	46000 UD	46000
4,6-Dinitro-2-methylphenol	46000 UD	46000

HPN: 2757

(continued)

Client ID: S2
Matrix: SOIL

Lab ID (HSN): 15277 DL Filename: 4067L16

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	18000 UD	18000
4-Bromophenyl-phenylether	18000 UD	18000
Hexachlorobenzene	18000 UD	18000
Pentachlorophenol	46000 UD	46000
Phenanthrene	18000 UD	18000
Anthracene	18000 UD	18000
Carbazole	18000 UD	18000
Di-n-butylphthalate	3000 JDB	18000
Fluoranthene	18000 UD	18000
Pyrene	2800 JD	18000
Butylbenzylphthalate	18000 UD	18000
3,3 ⁷ -Dichlorobenzidine	18000 UD	18000
Benz(a)anthracene	18000 UD	18000
Chrysene	1800 JD	18000
bis(2-Ethylhexyl)phthalate	2800 JD	18000
Di-n-octylphthalate	18000 UD	18000
Benzo(b) fluoranthene	2100 JD	18000
Benzo(k) fluoranthene	18000 UD	18000
Benzo(a)pyrene	18000 UD	18000
Indeno(1,2,3-cd)pyrene	18000 UD	18000
Dibenz(a,h)anthracene	18000 UD	18000
Benzo(g,h,i)perylene	18000 UD	18000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	49 % JD	25-121%
Phenol-d5	56 % JD	24-113%
2-Chlorophenol-d4	53 % JD	20-130%
1,2-Dichlorobenzene-d4	56 % JD	20-130%
Nitrobenzene-d5	56 % JD	23-120%
2-Fluorobiphenyl	66 % JD	30-115%
2,4,6-Tribromophenol	36 % JD	19-122*
Terphenyl-d14	90 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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Client ID: S3 Lab ID (HSN): 15278
Matrix: SOIL Filename: 4068K10

Date Sampled: 02/22/94 Sample Size: 15 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 1

Date Analyzed: 03/09/94 GPC Factor: 2 % Moisture: 30.97

Compounds:	ug/Kg (PPB)	EQL
Phenol	970 Ŭ	970
bis(2-Chloroethyl)ether	970 Ŭ	970
2-Chlorophenol	970 U	970
1,3-Dichlorobenzene	9 7 0 Ŭ	970
1,4-Dichlorobenzene	9 7 0 Ŭ	970
1,2-Dichlorobenzene	9 7 0 Ŭ	970
2-Methylphenol	970 Ŭ	970
2,2'-oxybis(1-Chloropropane)	970 Ŭ	970
4-Methylphenol	970 Ŭ	970
N-Nitroso-di-n-propylamine	970 U	970
Hexachloroethane	970 U	970
Nitrobenzene	970 U	970
Isophorone	970 U	970
2-Nitrophenol	970 U	970
2,4-Dimethylphenol	9 7 0 Ŭ	970
bis(2-Chloroethoxy)methane	970 U	970
2,4-Dichlorophenol	970 U	970
1,2,4-Trichlorobenzene	970 U	970
Naphthalene	9 7 0 Ŭ	970
4-Chloroaniline	. 970 บั	970
Hexachlorobutadiene	970 U	970
4-Chloro-3-methylphenol	970 Ŭ	970
2-Methylnaphthalene	970 U	970
Hexachlorocyclopentadiene	970 U	970
2,4,6-Trichlorophenol	9 7 0 Ŭ	970
2,4,5-Trichlorophenol	2400 U	2400
2-Chloronaphthalene	970 U	970
2-Nitroaniline	2400 U	2400
Dimethylphthalate	970 U	970
Acenaphthylene	970 Ŭ	970
2,6-Dinitrotoluene	970 U	970
3-Nitroaniline	2400 U	2400
Acenaphth ene	970 U	970
2,4-Dinitrophenol	2400 U	2400
4-Nitrophenol	2400 U	2400
Dibenzofuran	970 U	970
2,4-Dinitrotoluene	970 U	970
Diethylphthalate	970 Ŭ	970
4-Chlorophenyl-phenylether	970 U	970
Fluorene	970 U	970
4-Nitroaniline	2400 U	2400
4,6-Dinitro-2-methylphenol	2400 U	2400

(continued)

Huntingdon

Client ID: S3
Matrix: SOIL

Lab ID (HSN): 15278 Filename: 4068K10

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	970 U	970
4-Bromophenyl-phenylether	970 Ŭ	970
Hexachlorobenzene	970 Ŭ	970
Pentachlorophenol	2400 U	2400
Phenanthrene	880 J	970
Anthracene	140 J	970
Carbazole	970 Ŭ	970
Di-n-butylphthalate	1100 B	970
Fluoranthene	1700	970
Pyrene	3500 T Y	970
Butylbenzylphthalate	970 U Y	
3,3'-Dichlorobenzidine	970 U Y	970
Benz(a)anthracene	2000 T Y	970
Chrysene	2200 T Y	970
bis(2-Ethylhexyl)phthalate	3300 T Y	970
Di-n-octylphthalate	970 บ	970
Benzo(b) fluoranthene	3600	970
Benzo(k) fluoranthene	1100	970
Benzo(a) pyrene	2600	970
Indeno(1,2,3-cd)pyrene	2600	970
Dibenz(a,h)anthracene	970 U	970
Benzo(g,h,i)perylene	2600	970

Surrogate Recovery		QC LIMITS
2-Fluorophenol	25%	25-121%
Phenol-d5	34%	24-113%
2-Chlorophenol-d4	38%	20-130%
1,2-Dichlorobenzene-d4	40%	20-130%
Nitrobenzene-d5	29%	23-120%
2-Fluorobiphenyl	64%	30-115%
2,4,6-Tribromophenol	41%	19-122%
Terphenyl-d14	84*	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

Huntingdon_

Client ID: S3 Lab ID (HSN): 15278 DL Matrix: SOIL Filename: 4067L17

Sample Size: 15 grams Date Sampled: 02/22/94 Date Received: 02/24/94
Date Extracted: 03/03/94 500 uL Extract Vol.: Dil. Factor: 10 Date Analyzed: 03/09/94

GPC Factor: 2 % Moisture: 30.97

		
Compounds:	ug/Kg (PPB)	EQL
Phenol	9700 UD	9700
bis(2-Chloroethyl)ether	9700 UD	9700
2-Chlorophenol	9700 UD	9700
1,3-Dichlorobenzene	9700 UD	9700
1,4-Dichlorobenzene	9700 UD	9700
1,2-Dichlorobenzene	9700 UD	9700
2-Methylphenol	9700 UD	9700
2,2'-oxybis(1-Chloropropane)	9700 UD	9700
4-Methylphenol	9700 UD	9700
N-Nitroso-di-n-propylamine	9700 UD	9700
Hexachloroethane	9700 UD	9700
Nitrobenzene	. 9700 UD	9700
Isophorone	9700 UD	9700
2-Nitrophenol	9700 UD	9700
2,4-Dimethylphenol	9700 UD	9700
bis (2-Chloroethoxy) methane	9700 UD	9700
2,4-Dichlorophenol	9700 UD	9700
1,2,4-Trichlorobenzene	9700 UD	9700
Naphthalene	9700 UD	9700
4-Chloroaniline	9700 UD	9700
Hexachlorobutadiene	9700 UD	9700
4-Chloro-3-methylphenol	9700 UD ·	9700
2-Methylnaphthalene	9700 UD	9700
Hexachlorocyclopentadiene	9700 UD	9700
2,4,6-Trichlorophenol	9700 UD	9700
2,4,5-Trichlorophenol	24000 UD	24000
2-Chloronaphthalene	9700 UD	9700
2-Nitroaniline	24000 UD	24000
Dimethylphthalate	9700 UD	9700
Acenaphthylene	9700 UD	9700
2,6-Dinitrotoluene	9700 UD	9700
3-Nitroaniline	24000 UD	24000
Acenaphthene	9700 UD	9700
2,4-Dinitrophenol	24000 UD	24000
4-Nitrophenol	24000 UD	24000
Dibenzofuran	9700 UD	9700
2,4-Dinitrotoluene	9700 UD	9700
Diethylphthalate	9700 UD	9700
4-Chlorophenyl-phenylether	9700 UD	9700
Fluorene	9700 UD	9700
4-Nitroaniline	24000 UD	24000
4,6-Dinitro-2-methylphenol	24000 UD	24000
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Huntingdon

Client ID: S3
Matrix: SOIL

Dibenz(a,h)anthracene

Benzo(g,h,i)perylene

Lab ID (HSN): 15278 DL Filename: 4067L17

Jompounds:	ug/Kg (PP	PB)	EQL
	-3/3 (-,	202
N-Nitrosodiphenylamine	9700	യ	9700
4-Bromophenyl-phenylether	9700	UD	9700
Hexachlorobenzene	9700		9700
Pentachlorophenol	24000		24000
Phenanthrene		JD	9700
Anthracene	9700		9700
Carbazole	9700		9700
Di-n-butylphthalate	1600	JDB	9700
Fluoranthene	2700		9700
Pyrene	4100	JD	9700
Butylbenzylphthalate	9700	UD	9700
3,3'-Dichlorobenzidine	9700	UD	9700
Benz(a)anthracene	2300	JD	9700
Chrysene	2400	ற	9700
bis(2-Ethylhexyl)phthalate	6200	JD	9700
Di-n-octylphthalate	9700	UD	9700
Benzo(b) fluoranthene	4200	JD	9700
Benzo(k) fluoranthene	1300	JD	9700
Benzo(a)pyrene	2800	JD	9700
Indeno(1,2,3-cd)pyrene	3000	JD	9700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	36 % JD	25-121%
Phenol-d5	50 % JD	24-113%
2-Chlorophenol-d4	46 % JD	20-130%
1,2-Dichlorobenzene-d4	41 % JD	20-130%
Nitrobenzene-d5	46 % JD	23-120%
2-Fluorobiphenyl	75 %J D	30-115%
2,4,6-Tribromophenol	51 % JD	19-122%
Terphenyl-d14	110 % JD	18-137%

- TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)
- EQL = Estimated Quantitation Limit (lower calibration limit)
- U = Undetected at the given EQL
- J = Detected below the EQL (estimated value)
- E = Exceeds the upper calibration limit (estimated value)
- B = Also detected in the associated Blank
- D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

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HPN: 2757

9700 UD

2900 JD

9700

9700

Lab ID (HSN): 15279 Client ID: S4 Filename: 4068K07 Matrix: SOIL

Sample Size: 15.2 grams
Extract Vol.: 1500 uL
Dil. Factor: 1
GPC Factor: 2 Date Sampled: 02/22/94 Date Received: 02/24/94 Date Extracted: 03/03/94

Date Analyzed: 03/09/94 % Moisture: 5.88

Compounds:	ug/Kg	(PP	PB)	EQL
Phenol	21	00	U	2100
bis(2-Chloroethyl)ether	21	00	U	2100
2-Chlorophenol	21	00	U	2100
1,3-Dichlorobenzene	21	00	U	2100
1,4-Dichlorobenzene	21	00	U	2100
1,2-Dichlorobenzene	21	00	U	2100
2-Methylphenol	21	00	U	2100
2,2'-oxybis(1-Chloropropane)	21	00	บ	2100
4-Methylphenol	21	00	U	2100
N-Nitroso-di-n-propylamine	21	00	U	2100
Hexachloroethane	21	00	U	2100
Nitrobenzene	21	00	U	2100
Isophorone	21	00	U	2100
2-Nitrophenol	21	00	U	2100
2,4-Dimethylphenol	21	00	U	2100
bis(2-Chloroethoxy)methane	21	00	U	2100
2,4-Dichlorophenol	21	00	U	2100
1,2,4-Trichlorobenzene	21	00	U	2100
Naphthalene	21	00	U	2100
4-Chloroaniline	21	00	U	2100
Hexachlorobutadiene		00		2100
4-Chloro-3-methylphenol		00		2100
2-Methylnaphthalene		00		2100
Hexachlorocyclopentadiene		.00		2100
2,4,6-Trichlorophenol		00		2100
2,4,5-Trichlorophenol		00		5200
2-Chloronaphthalene		00		2100
2-Nitroaniline		00		5200
Dimethylphthalate		.00		2100
Acenaphthylene		00		2100
2,6-Dinitrotoluene		.00		2100
3-Nitroaniline		00		5200
Acenaphthene		00		2100
2,4-Dinitrophenol		00		5200
4-Nitrophenol		00		5200
Dibenzofuran		00		2100
2,4-Dinitrotoluene		.00		2100
Diethylphthalate		.00		2100
4-Chlorophenyl-phenylether		.00		2100
Fluorene		.00		2100
4-Nitroaniline		00		5200
4,6-Dinitro-2-methylphenol	52	00	J	5200

(continued)

Huntingdon

Client ID: S4
Matrix: SOIL

Lab ID (HSN): 15279 Filename: 4068K07

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	2100 U	2100
4-Bromophenyl-phenylether	2100 U	2100
Hexachlorobenzene	2100 U	2100
Pentachlorophenol	5200 U	5200
Phenanthrene	1800 J	2100
Anthracene	240 J	2100
Carbazole	2100 U	2100
Di-n-butylphthalate	2100 U	2100
Fluoranthene	2100 U	
Pyrene	2100 U Y	
Butylbenzylphthalate	2100 U Y	
3,3'-Dichlorobenzidine	2100 U Y	
Benz(a)anthracene	2100 U Y	
Chrysene	2100 U Y	
bis(2-Ethylhexyl)phthalate	6300 T Y	
Di-n-octylphthalate	2100 U	2100
Benzo(b) fluoranthene	4200	2100
Benzo(k) fluoranthene	1300 Ј	2100
Benzo(a) pyrene	2500	2100
Indeno(1,2,3-cd)pyrene	2600	2100
Dibenz(a,h)anthracene	2100 U	2100
Benzo(g,h,i)perylene	3100	2100

Surrogate Recovery		QC LIMITS
2-Fluorophenol	2 4 *J	25-121%
Phenol-d5	48%	24-113*
2-Chlorophenol-d4	46*	20-130%
1,2-Dichlorobenzene-d4	50 % J	20-130%
Nitrobenzene-d5	42 % J	23-120%
2-Fluorobiphenyl	107%	30-115%
2,4,6-Tribromophenol	52%	19-122%
Terphenyl-d14	115%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

/ = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

Huntingdon HPN: 2757

Client ID: S4 Lab ID (HSN): 15279 DL Matrix: SOIL Filename: 4067L14

Date Sampled: 02/22/94 Sample Size. 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1500 uL

Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQL
Phenol bis(2-Chloroethyl)ether 2-Chlorophenol 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 2-Methylphenol 2,2'-oxybis(1-Chloropropane) 4-Methylphenol N-Nitroso-di-n-propylamine Hexachloroethane Nitrobenzene Isophorone 2-Nitrophenol 2,4-Dimethylphenol bis(2-Chloroethoxy)methane 2,4-Dichlorophenol 1,2,4-Trichlorobenzene Naphthalene 4-Chloroaniline Hexachlorobutadiene 4-Chloro-3-methylphenol 2-Methylnaphthalene Hexachlorocyclopentadiene 2,4,6-Trichlorophenol 2,4,5-Trichlorophenol 2-Chloronaphthalene 2-Nitroaniline Dimethylphthalate	21000 UD 21000 UD	21000 21000
Acenaphthylene 2,6-Dinitrotoluene 3-Nitroaniline Acenaphthene	21000 UD 21000 UD 21000 UD 52000 UD 21000 UD	21000 21000 52000 21000
2,4-Dinitrophenol 4-Nitrophenol Dibenzofuran 2,4-Dinitrotoluene Diethylphthalate 4-Chlorophenyl-phenylether Fluorene 4-Nitroaniline	52000 UD 52000 UD 21000 UD 21000 UD 21000 UD 21000 UD 21000 UD 52000 UD	52000 52000 21000 21000 21000 21000 21000 52000
4,6-Dinitro-2-methylphenol	52000 UD	52000

(continued)

Client ID: S4
Matrix: SOIL

Lab ID (HSN): 15279 DL Filename: 4067L14

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	21000 UD	21000
4-Bromophenyl-phenylether	21000 UD	21000
Hexachlorobenzene	21000 UD	21000
Pentachlorophenol	52000 UD	52000
Phenanthrene	2400 JD	21000
Anthracene	21000 UD	21000
Carbazole	21000 UD	21000
Di-n-butylphthalate	21000 UD	21000
Fluoranthene	3800 JD	21000
Pyrene	7300 JD	21000
Butylbenzylphthalate	21000 UD	21000
3,3'-Dichlorobenzidine	21000 UD	21000
Benz (a) anthracene	2500 JD	21000
Chrysene	2800 JD	21000
bis(2-Ethylhexyl)phthalate	14000 JD	21000
Di-n-octylphthalate	21000 UD	21000
Benzo(b) fluoranthene	· 4000 JD	21000
Benzo(k) fluoranthene	21000 UD	21000
Benzo(a) pyrene	2400 JD	21000
Indeno(1,2,3-cd)pyrene	21000 UD	21000
Dibenz(a,h) anthracene	21000 UD	21000
Benzo(g,h,i)perylene	21000 UD	21000
	21000 02	21000

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl 2,4,6-Tribromophenol	28%JD 62%JD 52%JD 54%JD 57%JD 102%JD 54%JD	QC LIMITS 25-121% 24-113% 20-130% 20-130% 23-120% 30-115% 19-122%
Terphenyl-d14	126*JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

Huntingdon

Client ID: S4 MS Lab ID (HSN): 15279MS
Matrix: S0IL Filename: 4068K05

Date Sampled: 02/22/94 Sample Size: 15.6 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL
Date Extracted: 03/03/94 Dil. Factor: 1

Date Analyzed: 03/09/94 GPC Factor: 2 % Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQL
•		
Phenol	180 J	1400
bis(2-Chloroethyl)ether	1400 U	1400
2-Chlorophenol	1400 U	1400
1,3-Dichlorobenzene	1400 U	1400
1,4-Dichlorobenzene	240 J	1400
1,2-Dichlorobenzene	1400 U	1400
2-Methylphenol	1400 U	1400
2,2'-oxybis(1-Chloropropane)	1400 U	1400
4-Methylphenol	1400 U	1400
N-Nitroso-di-n-propylamine	290 J	1400
Hexachloroethane	1400 U	1400
Nitrobenzene	1400 U	1400
Isophorone	1400 U	1400
2-Nitrophenol	1400 U	1400
2,4-Dimethylphenol	1400 U	1400
bis(2-Chloroethoxy)methane	1400 U	1400
2,4-Dichlorophenol	1400 U	1400
1,2,4-Trichlorobenzene	290 J	1400
Naphthalene	1400 U	1400
4-Chloroaniline	1400 U	1400
Hexachlorobutadiene	1400 U	1400
4-Chloro-3-methylphenol	1400 U	1400
2-Methylnaphthalene	1400 U	1400
Hexachlorocyclopentadiene	1400 U	1400
2,4,6-Trichlorophenol	1400 U	1400
2,4,5-Trichlorophenol	3400 U	3400
2-Chloronaphthalene	1400 U	1400
2-Nitroaniline	3400 U	3400
Dimethylphthalate	1400 U	1400
Acenaphthylene	1400 U	1400
2,6-Dinitrotoluene	1400 U	1400
3-Nitroaniline	3400 U	3400
Acenaphth ene	3 4 0 J	1400
2,4-Dinitrophenol	3400 U	3400
4-Nitrophenol	380 J	3400
Dibenzofuran	1400 U	1400
2,4-Dinitrotoluene	320 J	1400
Diethylphthalate	1400 U	1400
4-Chlorophenyl-phenylether	1400 U	1400
Fluorene	1400 U	1400
4-Nitroaniline	3400 U	3400
4,6-Dinitro-2-methylphenol	3400 U	3400
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Huntingdon

Client ID: S4 MS Lab ID (HSN): 15279MS Matrix: SOIL Filename: 4068K05

Compounds:	ug/Kg (PF	PB) EQL
N-Nitrosodiphenylamine	1400	U 1400
4-Bromophenyl-phenylether	1400	U 1400
Hexachlorobenzene	1400	U 1400
Pentachlorophenol	3400	U 3400
Phenanthrene	650	J 1400
Anthracene	150	J 1400
Carbazole	1400	U 1400
Di-n-butylphthalate	1400	U 1400
Fluoranthene	640	J 1400
Pyrene	1300	J 1400
Butylbenzylphthalate	1400	U 1400
3,3'-Dichlorobenzidine	1400	U 1400
Benz(a)anthracene	370	J 1400
Chrysene	400	J 1400
bis(2-Ethylhexyl)phthalate	1200	J 1400
Di-n-octylphthalate	1400	U 1400
Benzo(b)fluoranthene	480	
Benzo(k)fluoranthene	160	
Benzo(a)pyrene	310	
Indeno(1,2,3-cd)pyrene	330	
Dibenz(a,h)anthracene	1400	
Benzo(g,h,i)perylene	1400	U 1400

Surrogate Recovery		QC LIMITS
2-Fluorophenol	1 % J	25-121
Phenol-d5	3 % J	24-113%
2-Chlorophenol-d4	3 % J	20-130%
1,2-Dichlorobenzene-d4	7 % J	20-130%
Nitrobenzene-d5	8 % J	23-120%
2-Fluorobiphenyl	9 % J	30-115%
2,4,6-Tribromophenol	0 % U	19-122%
Terphenyl-d14	10 % J	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Client ID: S4 MSD Lab ID (HSN): 15279MSD Matrix: SOIL Filename: 4068K06

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 2000 uL
Date Extracted: 03/03/94 Dil. Factor: 1

Date Analyzed: 03/09/94 GFC Factor: 2 % Moisture: 5.88

Compounds:	ug/Kg (PPB)	EQL
Phenol	6100	2800
bis(2-Chloroethyl)ether	2800 U	2800
2-Chlorophenol	2500 J	2800
1,3-Dichlorobenzene	2800 U	2800
1,4-Dichlorobenzene	1700 J	2800
1,2-Dichlorobenzene	2800 U	2800
2-Methylphenol	2800 U	2800
2,2'-oxybis(1-Chloropropane)	2800 U	2800
4-Methylphenol	2800 U	2800
N-Nitroso-di-n-propylamine	2100 J	2800
Hexachloroethane	2800 U	2800
Nitrobenzene	2800 U	2800
Isophorone	2800 U	2800
2-Nitrophenol	2800 U	2800
2,4-Dimethylphenol	2800 U	2800
bis(2-Chloroethoxy)methane	2800 U	2800
2,4-Dichlorophenol	2800 U	2800
1,2,4-Trichlorobenzene	2800	2800
Naphthalene	2800 U	2800
4-Chloroaniline	2800 U	2800
Hexachlorobutadiene	2800 U	2800
4-Chloro-3-methylphenol	2900	2800
2-Methylnaphthalene	2800 U	2800
Hexachlorocyclopentadiene	2800 U	2800
2,4,6-Trichlorophenol	2800 U	2800
2,4,5-Trichlorophenol	6900 U	6900
2-Chloronaphthalene	2800 U	2800
2-Nitroaniline	6900 U	6900
Dimethylphthalate	2800 U	2800
Acenaphthylene	2800 U	2800
2,6-Dinitrotoluene	2800 U	2800
3-Nitroaniline	6900 U	6900 2800
Acenaphthene	2900 6900 U	6900
2,4-Dinitrophenol	6900 U	6900
4-Nitrophenol	2800 U	2800
Dibenzofuran	3100	2800
2,4-Dinitrotoluene	2800 U	2800
Diethylphthalate 4-Chlorophenyl-phenylether	2800 U	2800
fluorene	2800 U	2800
4-Nitroaniline	6900 U	6900
4,6-Dinitro-2-methylphenol	6900 U	6900
4,0-Difficio-7-Weeth ibuenot		~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~

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Huntingdon

Client ID: S4 MSD Matrix: SOIL

Lab ID (HSN): 15279MSD Filename: 4068K06

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	2800 U	2800
4-Bromophenyl-phenylether	2800 U	2800
Hexachlorobenzene	2800 U	2800
Pentachlorophenol	6900 U	6900
Phenanthrene	1600 J	2800
Anthracene	2800 Ŭ	2800
Carbazole	2800 U	2800
Di-n-butylphthalate	17000 B	2800
Fluoranthene	2800 U	2800
Pyrene	6900 J Y	2800
Butylbenzylphthalate	2800 U Y	
3,3'-Dichlorobenzidine	2800 U Y	
Benz(a)anthracene	2400 J Y	
Chrysene	3000 1 Y	
bis(2-Ethylhexyl)phthalate	6500 T Y	2800
Di-n-octylphthalate	2800 U	2800
Benzo(b)fluoranthene	4800	2800
Benzo(k)fluoranthene	1600 J	2800
Benzo(a)pyrene	2600 J	2800
Indeno(1,2,3-cd)pyrene	2800	2800
Dibenz(a,h)anthracene	2800 U	2800
Benzo(g,h,i)perylene	2900	2800
	· ·	

Surrogate Recovery		QC LIMITS
2-Fluorophenol	33 % J	25-121%
Phenol-d5	52 % J	24-113%
2-Chlorophenol-d4	48 % J	20-130%
1,2-Dichlorobenzene-d4	5 4% J	20-130%
Nitrobenzene-d5	58 % J	23-120%
2-Fluorobiphenyl	94*	30-115%
2,4,6-Tribromophenol	38 % J	19-122%
Terphenyl-d14	77 % J	Y 18-137*

- TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)
- EQL = Estimated Quantitation Limit (lower calibration limit)
- U = Undetected at the given EQL
- J = Detected below the EQL (estimated value)
- E = Exceeds the upper calibration limit (estimated value)
- B = Also detected in the associated Blank
- ? = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,
November 1986, 3rd Edit::n.

Huntingdon

Client ID: S5 Lab ID (HSN): 15283 Matrix: SOIL Filename: 4068K11

Sample Size: 15.4 grams
Extract Vol.: 1000 uL
Dil. Factor: 1 Date Sampled: 02/22/94 Date Received: 02/24/94
Date Extracted: 03/03/94

Date Analyzed: 03/09/94 GPC Factor: 2

GPC Factor: 2
% Moisture: 22.42

Compounds:	ug/Kg (PPB)	EQL
Phenol	190 J	1700
bis(2-Chloroethyl)ether	1700 U	1700
2-Chlorophenol	1700 U	1700
1,3-Dichlorobenzene	1700 U	1700
1,4-Dichlorobenzene	1700 U	1700
1,2-Dichlorobenzene	1700 U	1700
2-Methylphenol	1700 U	1700
2,2'-oxybis(1-Chloropropane)	1700 U	1700
4-Methylphenol	1700 U	1700
N-Nitroso-di-n-propylamine	1700 U	1700
Hexachloroethane	1700 U	1700
Nitrobenzene	. 1700 U	1700
Isophorone	1700 U	1700
2-Nitrophenol	1700 U	1700
2,4-Dimethylphenol	1700 U	1700
bis(2-Chloroethoxy) methane	1700 U	1700
2,4-Dichlorophenol	1700 U	1700
1,2,4-Trichlorobenzene	720 J	1700
Naphthalene	220 J	1700
4-Chloroaniline	1700 U	1700
Hexachlorobutadiene	1700 U	1700
4-Chloro-3-methylphenol	1700 U	1700
2-Methylnaphthalene	1700 U	1700
Hexachlorocyclopentadiene	1700 U	1700
2,4,6-Trichlorophenol	1700 U	1700
2,4,5-Trichlorophenol	4200 U	4200
2-Chloronaphthalene	1700 U	1700
2-Nitroaniline	4200 U	4200
Dimethylphthalate	1700 U	1700 1700
Acenaphthylene	1700 U 1700 U	1700
2,6-Dinitrotoluene	4200 U	4200
3-Nitroaniline	340 J	1700
Acenaphthene	4200 U	4200
2,4-Dinitrophenol	4200 U	4200
4-Nitrophenol	1700 U	1700
Dibenzofuran	1700 U	1700
2,4-Dinitrotoluene	1700 U	1700
Diethylphthalate	1700 U	1700
4-Chlorophenyl-phenylether	410 J	1700
Fluorene	4200 U	4200
4-Nitroaniline	4200 U	4200
4,6-Dinitro-2-methylphenol	7200 0	4200

(continued)

Client ID: S5
Matrix: SOIL

Lab ID (HSN): 15283 Filename: 4068K11

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1700 U	1700
4-Bromophenyl-phenylether	1700 U	1700
Hexachlorobenzene	1700 U	1700
Pentachlorophenol	4200 U	4200
Phenanthrene	3500	1700
Anthracene	830 J	1700
Carbazole	1700 U	1700
Di-n-butylphthalate	1300 ЈВ	1700
Fluoranthene	3900	1700
Pyrene	7900 = Y	1700
Butylbenzylphthalate	1700 U Y	1700
3,3'-Dichlorobenzidine	1700 U Y	1700
Benz(a)anthracene	3800 ~ Y	1700
Chrysene	3800 🕻 Y	
bis(2-Ethylhexyl)phthalate	30000 E Y	
Di-n-octylphthalate	1700 U	1700
Benzo(b) fluoranthene	5800	1700
Benzo(k) fluoranthene	2200	1700
Benzo(a) pyrene	4400	1700
Indeno(1,2,3-cd)pyrene	4200	1700
Dibenz(a,h)anthracene	1700 U	1700
Benzo(g,h,i)perylene	4300	1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30∜	25-121%
Phenol-d5	37%	24-113%
2-Chlorophenol-d4	40%	20-130%
1,2-Dichlorobenzene-d4	46%	20-130%
Nitrobenzene-d5	31 % J	23-120%
2-Fluorobiphenyl	72%	30-115*
2,4,6-Tribromophenol	32*	19-122%
Terphenyl-d14	77%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y = Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Eva. Lating Solid Waste", SW-846,

November 1986, 3rd Edition

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Client ID: S5 Lab ID (HSN): 15283 DL Filename: 4067L18

Matrix: SOIL Date Sampled: 02/22/94 Sample Size: 15.4 grams Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Date Analyzed: 03/09/94 Dil. Factor: 10 GPC Factor: 2 % Moisture: 22.42

Compounds:	ug/Kg (PPB)	EQL
Phenol	17000 UD	17000
bis(2-Chloroethyl)ether	17000 UD	17000
2-Chlorophenol	17000 UD	17000
1,3-Dichlorobenzene	17000 UD	17000
1,4-Dichlorobenzene	17000 UD	17000
1,2-Dichlorobenzene	17000 UD	17000
2-Methylphenol	17000 UD	17000
2,2'-oxybis(1-Chloropropane)	17000 UD	17000
4-Methylphenol	17000 UD	17000
N-Nitroso-di-n-propylamine	17000 UD	17000
Hexachloroethane	17000 UD	17000
Nitrobenzene	17000 UD	17000
Isophorone	17000 UD	17000
2-Nitrophenol	17000 UD	17000
2,4-Dimethylphenol	17000 UD	17000
bis(2-Chloroethoxy)methane	17000 UD	17000
2,4-Dichlorophenol	17000 UD	17000
1,2,4-Trichlorobenzene	17000 UD	17000
Naphthalene	17000 UD	17000
4-Chloroaniline	17000 UD	17000
Hexachlorobutadiene	17000 UD	17000
4-Chloro-3-methylphenol	17000 UD	17000
2-Methylnaphthalene	17000 UD	17000
Hexachlorocyclopentadiene	17000 UD	17000
2,4,6-Trichlorophenol	17000 UD	17000
2,4,5-Trichlorophenol	42000 UD	42000
2-Chloronaphthalene	17000 UD	17000
2-Nitroaniline	42000 UD	42000
Dimethylphthalate	17000 UD	17000
Acenaphthylene	17000 UD	17000
2,6-Dinitrotoluene	17000 UD	17000
3-Nitroaniline	42000 UD	42000
Acenaphthen e	17000 UD	17000
2,4-Dinitrophenol	42000 UD	42000
4-Nitrophenol	42000 UD	42000
Dibenzofuran	17000 UD	17000
2,4-Dinitrotoluene	17000 UD	17000
Diethylphthalate	17000 UD	17000
4-Chlorophenyl-phenylether	17000 UD	17000
Fluorene	17000 UD	17000
4-Nitroaniline	42000 UD	42000
4,6-Dinitro-2-methylphenol	42000 UD	42000

(continued)

Client ID: S5 Lab ID (HSN): 15283 DL Matrix: SOIL Filename: 4067L18

Compounds:	ug/Kg (PE	B)	EQL
N-Nitrosodiphenylamine	17000	UD	17000
4-Bromophenyl-phenylether	17000	UD	17000
Hexachlorobenzene	17000	Œ	17000
Pentachlorophenol	42000	Œ	42000
Phenanthrene	4500	JD	17000
Anthracene	17000	S	17000
Carbazole	17000	UD	17000
Di-n-butylphthalate	2800	JDB	17000
Fluoranthene	5800	JD	17000
Pyrene	10000	JD	17000
Butylbenzylphthalate	17000	യ	17000
3,3'-Dichlorobenzidine	17000	UD	17000
Benz(a)anthracene	4100	JD	17000
Chrysene	3800	JD	17000
bis(2-Ethylhexyl)phthalate	56000	D	17000
Di-n-octylphthalate	17000	W	17000
Benzo(b)fluoranthene	6900	JD	17000
Benzo(k)fluoranthene	17000	Œ	17000
Benzo(a)pyrene	4500		17000
Indeno(1,2,3-cd)pyrene	5000		17000
Dibenz(a,h)anthracene	17000	UD	17000
Benzo(g,h,i)perylene	5000	JD	17000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	41%JD	25-121%
Phenol-d5	49 % JD	24-113%
2-Chlorophenol-d4	47 % JD	20-130%
1,2-Dichlorobenzene-d4	46 % JD	20-130%
Nitrobenzene-d5	48 % JD	23-120%
2-Fluorobiphenyl	66 % JD	30-115%
2,4,6-Tribromophenol	36 % JD	19-122%
Terphenyl-d14	96 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

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EQL = Estimated Quantitation Limit (lower calibration limit)

U - Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

⁻ Analysis at a secondary Dilution factor

Lab ID (HSN): 15287 Client ID: S6 Filename: 4068K12 Matrix: SOIL

Sample Size: 15.2 grams Date Sampled: 02/22/94 Extract Vol.: Date Received: 02/24/94 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 1 GPC Factor: Date Analyzed: 03/10/94 2

% Moisture: 17.96

Compounds:	ug/Kg (PPB)	EQ
Phenol	1600 U	160
bis(2-Chloroethyl)ether	1600 U	160
2-Chlorophenol	1600 U	160
1,3-Dichlorobenzene	1600 U	160
1,4-Dichlorobenzene	1600 U	160
1,2-Dichlorobenzene	1600 U	160
2-Methylphenol	1600 U	160
2,2'-oxybis(1-Chloropropane)	1600 U	160
4-Methylphenol	1600 U	160
N-Nitroso-di-n-propylamine	1600 U	160
Hexachloroethane	1600 U	160
Nitrobenzene	1600 U	160
Isophorone	1600 U	160
2-Nitrophenol	1600 U	160
2,4-Dimethylphenol	1600 U	160
bis (2-Chloroethoxy) methane	1600 U	160
2,4-Dichlorophenol	1600 U	160
1,2,4-Trichlorobenzene	690 J	160
Naphthalene	. 510 J	160
4-Chloroaniline	1600 U	160
Hexachlorobutadiene	1600 U	160
4-Chloro-3-methylphenol	1600 U	160
2-Methylnaphthalene	330 J	160
Hexachlorocyclopentadiene	1600 U	160
2,4,6-Trichlorophenol	1600 U	160
2,4,5-Trichlorophenol	4000 U	400
2-Chloronaphthalene	1600 U	160
2-Nitroaniline	4000 U	400
Dimethylphthalate	1600 U	160
Acenaphthylene	1600 U	160
2,6-Dinitrotoluene	1600 U	160
3-Nitroaniline	4000 U	400
Acenaphthene	740 J	160
2,4-Dinitrophenol	4000 U	400
4-Nitrophenol	4000 U	400
Dibenzofuran	760 J	160
2,4-Dinitrotoluene	1600 U	160
Diethylphthalate	1600 U	160
4-Chlorophenyl-phenylether	1600 U	160
Fluorene	1100 J	160
4-Nitroaniline	4000 U	400
4,6-Dinitro-2-methylphenol	4000 U	400

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(continued)

Client ID: S6 Lab ID (HSN): 15287 Matrix: SOIL Filename: 4068K12

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	1600 U	1600
4-Bromophenyl-phenylether	1600 U	1600
Hexachlorobenzene	1600 U	1600
Pentachlorophenol	4000 U	4000
Phenanthrene	7900	1600
Anthracene	2100	1600
Carbazole	760 J	1600
Di-n-butylphthalate	820 JB	1600
Fluoranthene	6700	1600
Pyrene	13000 E Y	1600
Butylbenzylphthalate	1600 U Y	1600
3,3'-Dichlorobenzidine	1600 U Y	1600
Benz(a)anthracene	6300 7 Y	1600
Chrysene	5700 🗸 Y	1600
bis(2-Ethylhexyl)phthalate	22000 E Y	1600
Di-n-octylphthalate	1600 U	1600
Benzo(b) fluoranthene	7100	1600
Benzo(k) fluoranthene	3000	1600
Benzo(a)pyrene	5700	1600
Indeno(1,2,3-cd)pyrene	4700	1600
Dibenz(a,h)anthracene	1600 U	1600
Benzo(g,h,i)perylene	4000	1600

Surrogate Recovery		QC LIMITS
2-Fluorophenol	45*	25-121%
Phenol-d5	47%	24-113%
2-Chlorophenol-d4	54*	20-130%
1,2-Dichlorobenzene-d4	67 %	20-130%
Nitrobenzene-d5	43%	23-120%
2-Fluorobiphenyl	77%	30-115%
2,4,6-Tribromophenol	31%	19-122%
Terphenyl-d14	82*	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y - Associated internal standard failed method criteria

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

Huntingdon

Client ID: S6 Lab ID (HSN): 15287 DL Matrix: SOIL Filename: 4067L19

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 1000 ul
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

% Moisture: 17.96

Compounds:	ug/Kg (PPB)	EQL
Phenol	16000 UD	16000
bis(2-Chloroethyl)ether	16000 UD	16000
2-Chlorophenol	16000 UD	16000
1,3-Dichlorobenzene	16000 UD	16000
1,4-Dichlorobenzene	16000 UD	16000
1,2-Dichlorobenzene	16000 UD	16000
2-Methylphenol	16000 UD	16000
2,2'-oxybis(1-Chloropropane)	16000 UD	16000
4-Methylphenol	16000 UD	16000
N-Nitroso-di-n-propylamine	16000 UD	16000
Hexachloroethane	16000 UD	16000
Nitrobenzene	16000 UD	16000
Isophorone	16000 UD	16000
2-Nitrophenol	16000 UD	16000
2,4-Dimethylphenol	16000 UD	16000
bis(2-Chloroethoxy)methane	16000 UD	16000
2,4-Dichlorophenol	16000 UD	16000
1,2,4-Trichlorobenzene	16000 UD	16000
Naphthalene	16000 UD	16000
4-Chloroaniline	16000 UD	16000
Hexachlorobutadiene	16000 UD	16000
4-Chloro-3-methylphenol	16000 UD	16000
2-Methylnaphthalene	16000 UD	16000
Hexachlorocyclopentadiene	16000 UD	16000
2,4,6-Trichlorophenol	16000 UD	16000
2,4,5-Trichlorophenol	40000 UD	40000
2-Chloronaphthalene	16000 UD	16000
2-Nitroaniline	40000 UD	40000
Dimethylphthalate	16000 UD	16000
Acenaphthylene	16000 UD	16000
2,6-Dinitrotoluene	16000 UD	16000
3-Nitroaniline	40000 UD	40000
Acenaphthene	16000 UD	16000
2,4-Dinitrophenol	40000 UD	40000
4-Nitrophenol	40000 UD	40000
Dibenzofuran	16000 UD	16000
2,4-Dinitrotoluene	16000 UD	16000
Diethylphthalate	16000 VD	16000
4-Chlorophenyl-phenylether	16000 UD	16000
Fluorene	16000 UD	
4-Nitroaniline	40000 UD	
4,6-Dinitro-2-methylphenol	40000 UD	40000

(continued)

Huntingdon

Client ID: S6
Matrix: SOIL

lab ID (HSN): 15287 DL
 Filename: 4067L19

Compounds:	ug/Kg (PF	PB)	EQL
N-Nitrosodiphenylamine 4-Bromophenyl-phenylether Hexachlorobenzene Pentachlorophenol Phenanthrene Anthracene Carbazole Di-n-butylphthalate Fluoranthene Pyrene Butylbenzylphthalate 3,3'-Dichlorobenzidine Benz(a)anthracene Chrysene bis(2-Ethylhexyl)phthalate Di-n-octylphthalate Benzo(b)fluoranthene Benzo(k)fluoranthene	16000 16000 40000 9700 2900 16000 1800 17000 16000 6400 5900 42000 16000 8000 3500	9999999999999999	16000 16000 16000 16000 16000 16000 16000 16000 16000 16000 16000 16000 16000
Benzo(a)pyrene Indeno(1,2,3-cd)pyrene Dibenz(a,h)anthracene Benzo(g,h,i)perylene	6000 5900 16000 5400	JD	16000 16000 16000 16000

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl 2,4,6-Tribromophenol	63%JD 71%JD 65%JD 66%JD 76%JD 35%JD	QC LIMITS 25-121% 24-113% 20-130% 20-130% 23-120% 30-115% 19-122%
Terphenyl-d14	108 %J D	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a try weight basis.

Reference: "EPA Test Methods for Evaluting Solid Waste", SW-846,
November 1986, 3rd Edit.

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EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

⁼ Analysis at a secondary Dilution factor

Lab ID (HSN): 15289 Client ID: S7 Matrix: SOIL Filename: 4068Kl3

Sample Size: 15.3 grams Date Sampled: 02/22/94 Date Received: 02/24/94 Extract Vol.: 500 ũL

Date Extracted: 03/03/94 Date Analyzed: 03/10/94 Dil. Factor: 1 GPC Factor: 2
% Moisture: 30.67

Compounds:	ug/Kg (PPB)	EQL
Phenol	9 4 0 U	940
bis(2-Chloroethyl)ether	9 4 0 U	940
2-Chlorophenol	9 4 0 U	940
1,3-Dichlorobenzene	9 4 0 U	940
1,4-Dichlorobenzene	9 4 0 U	940
1,2-Dichlorobenzene	9 40 U	940
2-Methylphenol	9 4 0 U	940
2,2'-oxybis(1-Chloropropane)	9 4 0 Ŭ	940
4-Methylphenol	940 U	940
N-Nitroso-di-n-propylamine	9 4 0 U	940
Hexachloroethane	940 U	940
Nitrobenzene	9 4 0 U	940
Isophorone	940 U	940
2-Nitrophenol	9 40 U	940
2,4-Dimethylphenol	9 4 0 U	940
bis(2-Chloroethoxy)methane	9 4 0 U	940
2,4-Dichlorophenol	9 4 0 U	940
1,2,4-Trichlorobenzene	430 J	940
Naphthalene	260 J	940
4-Chloroaniline	9 4 0 U	940
Hexachlorobutadiene	940 U	940
4-Chloro-3-methylphenol	9 4 0 U .	940
2-Methylnaphthalene	550 J	940
Hexachlorocyclopentadiene	9 4 0 U	940
2,4,6-Trichlorophenol	940 U	940
2,4,5-Trichlorophenol	2400 U	2400
2-Chloronaphthalene	940 Ŭ	940
2-Nitroaniline	2400 U	2400
Dimethylphthalate	940 U	940
Acenaphthylene	940 U	940
2,6-Dinitrotoluene	940 U	940
3-Nitroaniline	2400 U	2400
Acenaphthene	250 J	940
2,4-Dinitrophenol	2400 U	2400
4-Nitrophenol	2400 U	2400
Dibenzofuran	280 J	940
2,4-Dinitrotoluene	940 U	940
Diethylphthalate	940 U	940
4-Chlorophenyl-phenylether	9 4 0 U	940
Fluorene	290 J	940
4-Nitroaniline	2400 U	2400
4,6-Dinitro-2-methylphenol	2400 U	2400
,		

(continued)

Client ID: S7
Matrix: SOIL

Lab ID (HSN): 15289 Filename: 4068K13

Compounds:	ug/Kq (PPB)	FOT
compounds.	dg/kg (PPB)	EQL
N-Nitrosodiphenylamine	9 4 0 U	940
4-Bromophenyl-phenylether	940 U	940
Hexachlorobenzene	940 U	940
Pentachlorophenol	2400 U	2400
Phenanthrene	2100	940
Anthracene	580 J	940
Carbazole	940 U	940
Di-n-butylphthalate	800 JB	940
Fluoranthene	2700	940
Pyrene	5400 T Y	940
Butylbenzylphthalate	940 U Y	940
3,3'-Dichlorobenzidine	940 U Y	940
Benz(a)anthracene	2100 - Y	940
Chrysene	2300 7 Y	940
bis(2-Ethylhexyl)phthalate	9400 E Y	940
Di-n-octylphthalate	9 4 0 U	940
Benzo(b) fluoranthene	3600	940
Benzo(k)fluoranthene	1200	940
Benzo(a)pyrene	2200	940
Indeno(1,2,3-cd)pyrene	2100	940
Dibenz(a,h)anthracene	460 J	940
Benzo(g,h,i)perylene	1900	940

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30%	25-121%
Phenol-d5	36 %	24-113%
2-Chlorophenol-d4	41%	20-130%
1,2-Dichlorobenzene-d4	41%	20-130%
Nitrobenzene-d5	30%	23-120%
2-Fluorobiphenyl	71%	30-115%
2,4,6-Tribromophenol	47%	19-122*
Terphenyl-d14	75%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLMO1)

Note: All results are reported on a try weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,
November 1986, 3rd Edit:

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EQL = Estimated Quantitation Limit | lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit estimated value).

B = Also detected in the associated Blank

^{! =} Associated internal standard fa:.ed method criteria

Client ID: S7 Lab ID (HSN): 15289 DL Matrix: SOIL Filename: 4067L20

Date Sampled: 02/22/94 Sample Size: 15.3 grams Date Received: 02/24/94 Extract Vol.: 500 ŭL

Date Extracted: 03/03/94 Dil. Factor: 10 GPC Factor: Date Analyzed: 03/09/94 2

% Moisture: 30.67

Compounds:	ug/Kg (P	PB)	EQL
Phenol	9400	מוז	9400
bis(2-Chloroethyl)ether	9400		9400
2-Chlorophenol	9400		9400
1,3-Dichlorobenzene	9400		9400
1,4-Dichlorobenzene	9400		9400
1,2-Dichlorobenzene	9400		9400
2-Methylphenol	9400		9400
2,2'-oxybis(1-Chloropropane)	9400		9400
4-Methylphenol	9400		9400
N-Nitroso-di-n-propylamine	9400		9400
Hexachloroethane	9400		9400
Nitrobenzene	9400	UD	9400
Isophorone	9400	UD	9400
2-Nitrophenol	9400		9400
2,4-Dimethylphenol	9400		9400
bis(2-Chloroethoxy)methane	9400		9400
2,4-Dichlorophenol	9400		9400
1,2,4-Trichlorobenzene	9400		9400
Naphthalene	9400		9400
4-Chloroaniline	9400		9400
Hexachlorobutadiene		ω ·	9400
4-Chloro-3-methylphenol	9400		9400
2-Methylnaphthalene	9400		9400
Hexachlorocyclopentadiene	9400		9400
2,4,6-Trichlorophenol	9400		9400
2,4,5-Trichlorophenol	24000	UD	24000
2-Chloronaphthalene	9400	UD	9400
2-Nitroaniline	24000	UD	24000
Dimethylphthalate	9400	UD	9400
Acenaphthylene	9400	UD	9400
2,6-Dinitrotoluene	9400	UD	9400
3-Nitroaniline	24000	UD	24000
Acenaphthene	9400	WD	9400
2,4-Dinitrophenol	24000		24000
4-Nitrophenol	24000	UD	24000
Dibenzofuran	9400	UD	9400
2,4-Dinitrotoluene	9400	UD	9400
Diethylphthalate	9400	UD	9400
4-Chlorophenyl-phenylether	9400	UD .	9400
Fluorene	9400	UD ·	9400
4-Nitroaniline	24000		24000
4,6-Dinitro-2-methylphenol	24000	UD.	24000

Client ID: S7 Lab ID (HSN): 15289 DL Matrix: SOIL Filename: 4067L20

Compounds:	ug/Kg (PP	B) EQL
N-Nitrosodiphenylamine	9400	UD 9400
4-Bromophenyl-phenylether	9400	UD 9400
Hexachlorobenzene	9400	
Pentachlorophenol	24000	
Phenanthrene	2200	
Anthracene	9400	
Carbazole	9400	
Di-n-butylphthalate	1100	
Fluoranthene	3200	
Pyrene		JD 9400
Butylbenzylphthalate	9400	
3,3'-Dichlorobenzidine	9400	
Benz(a)anthracene	2200	
Chrysene	2400	-
bis(2-Ethylhexyl)phthalate	18000	
Di-n-octylphthalate	9400	
Benzo(b) fluoranthene	3700	
Benzo(k) fluoranthene	940	
Benzo(a) pyrene	2400	
Indeno(1,2,3-cd)pyrene	2400	
Dibenz(a,h)anthracene	9400	
Benzo(g,h,i)perylene	2500	

Surrogate Recovery		QC LIMITS
2-Fluorophenol	39 ∜ JD	25-121%
Phenol-d5	51 % JD	24-113%
2-Chlorophenol-d4	46 % JD	20-130%
1,2-Dichlorobenzene-d4	39 ∜ JD	20-130%
Nitrobenzene-d5	45 % JD	23-120%
2-Fluorobiphenyl	70 %J D	30-115%
2,4,6-Tribromophenol	55 ≹J D	19-122%
Terphenyl-d14	108 % JD	18-137%

- TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)
- EQL = Estimated Quantitation Limit (lower calibration limit)
 - U = Undetected at the given EQL
 - J = Detected below the EQL (estimated value)
 - E = Exceeds the upper calibration limit (estimated value)
 - B = Also detected in the associated Blank
 - D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

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Client ID: S8 Lab ID (HSN): 15290 Matrix: SOIL Filename: 4068K14

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/10/94 GPC Factor: 2
% Moisture: 23.54

Compounds:	ug/Kg (PPB)	EQL
Phenol	1700 U	1700
bis(2-Chloroethyl)ether	1700 U	1700
2-Chlorophenol	1700 U	1700
1,3-Dichlorobenzene	1700 U	1700
1,4-Dichlorobenzene	1700 U	1700
1,2-Dichlorobenzene	1700 U	1700
2-Methylphenol	1700 U	1700
2,2'-oxybis(1-Chloropropane)	1700 U	1700
4-Methylphenol	1700 U	1700
N-Nitroso-di-n-propylamine	1700 U	1700
Hexachloroethane	1700 U	1700
Nitrobenzene	1700 U	1700
Isophorone	1700 U	1700
2-Nitrophenol	1700 U	1700
2,4-Dimethylphenol	1700 U	1700
bis(2-Chloroethoxy)methane	1700 U	1700
2,4-Dichlorophenol	1700 U	1700
1,2,4-Trichlorobenzene	1700 U	1700
Naphthalene	250 J	1700
4-Chloroaniline	· 1700 U	1700
Hexachlorobutadiene	1700 U	1700
4-Chloro-3-methylphenol	1700 U	1700
2-Methylnaphthalene	480 J	1700
Hexachlorocyclopentadiene	1700 U	1700
2,4,6-Trichlorophenol	1700 U	1700
2,4,5-Trichlorophenol	4300 U	4300
2-Chloronaphthalene	1700 U	1700
2-Nitroaniline	4300 U	4300
Dimethylphthalate	1700 U	1700
Acenaphthyl ene	1700 U	1700
2,6-Dinitrotoluene	1700 U	1700
3-Nitroaniline	4300 U	4300
Acenaphthen e	520 J	1700
2,4-Dinitrophenol	4300 U	4300
4-Nitrophenol	4300 U	4300
Dibenzofuran	230 J	1700
2,4-Dinitrotoluene	1700 U	1700
Diethylphthalate	1700 U	1700
4-Chlorophenyl-phenylether	1700 U	1700
Fluorene		. 1700
4-Nitroaniline	4300 U	4300
4,6-Dinitro-2-methylphenol	4300 U	4300

(continued)

Client ID: S8
Matrix: SOIL

Lab ID (HSN): 15290 Filename: 4068K14

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine 4-Bromophenyl-phenylether Hexachlorobenzene Pentachlorophenol Phenanthrene Anthracene Carbazole Di-n-butylphthalate Fluoranthene Pyrene Butylbenzylphthalate 3,3'-Dichlorobenzidine Benz(a) anthracene Chrysene bis(2-Ethylhexyl) phthalate Di-n-octylphthalate Benzo(b) fluoranthene Benzo(k) fluoranthene Benzo(a) pyrene Indeno(1,2,3-cd) pyrene Dibenz(a,h) anthracene	1700 U 1700 U 1700 U 1700 U 4300 U 2900 420 J 1700 U 3900 B 3600 7600 T Y 4500 T Y 4700 T Y 4700 T Y 4700 T Y 4700 T Y 4700 T Y 4700 U 7600 2400 5200 6300 1100 J	1700 1700 1700 4300 1700 1700 1700 1700 1700 1700 1700
Benzo(g,h,i)perylene	6100	1700

Surrogate Recovery		QC LIMITS
2-Fluorophenol	43%	25-121%
Phenol-d5	45%	24-113%
2-Chlorophenol-d4	55%	20-130%
1,2-Dichlorobenzene-d4	56 %	20-130%
Nitrobenzene-d5	36 % J	23-120%
2-Fluorobiphenyl	· 75 %	30-115%
2,4,6-Tribromophenol	61 %	19-122%
Terphenyl-d14	96*	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

TToutingdon

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

Y - Associated internal standard failed method criteria

Client ID: S8 Lab ID (HSN): 15290 DL Matrix: SOIL Filename: 4067L21

Date Sampled: 02/22/94 Sample Size: 15.3 grams
Date Received: 02/24/94 Extract Vol.: 1000 uL

Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 23.54

Compounds:	ug/Kg (PPB)	EQL
Phenol	17000 UD	17000
bis(2-Chloroethyl)ether	17000 UD	17000
2-Chlorophenol	17000 UD	17000
1,3-Dichlorobenzene	17000 UD	17000
1,4-Dichlorobenzene	17000 UD	17000
1,2-Dichlorobenzene	17000 UD	17000
2-Methylphenol	17000 UD	17000
2,2'-oxybis(1-Chloropropane)	17000 UD	17000
4-Methylphenol	17000 UD	17000
N-Nitroso-di-n-propylamine	17000 UD	17000
Hexachloroethane	17000 UD	17000
Nitrobenzene	17000 UD	17000
Isophorone	17000 UD	17000
2-Nitrophenol	17000 UD	17000
2,4-Dimethylphenol	17000 UD	17000
bis (2-Chloroethoxy) methane	17000 UD	17000
2,4-Dichlorophenol	17000 UD	17000
1,2,4-Trichlorobenzene	17000 UD	17000
Naphthalene	17000 UD	17000
4-Chloroaniline	17000 UD	17000
Hexachlorobutadiene	17000 UD	17000
4-Chloro-3-methylphenol	17000 UD	17000
2-Methylnaphthalene	17000 UD	17000
Hexachlorocyclopentadiene	17000 UD	17000
2,4,6-Trichlorophenol	17000 UD	17000
2,4,5-Trichlorophenol	43000 UD	43000
2-Chloronaphthalene	17000 UD	17000
2-Nitroaniline	43000 UD	43000
Dimethylphthalate	17000 UD	17000
Acenaphthylene	17000 UD	17000
2,6-Dinitrotoluene	17000 UD	17000
3-Nitroaniline	43000 UD	43000
Acenaphthene	17000 UD	17000
2,4-Dinitrophenol	43000 UD	43000
4-Nitrophenol	43000 UD	43000
Dibenzofuran	17000 UD	17000
2,4-Dinitrotoluene	17000 UD	17000
Diethylphthalate	17000 UD	17000
4-Chlorophenyl-phenylether	17000 UD	17000
Fluorene	17000 UD	17000
4-Nitroaniline	43000 UD	43000
4,6-Dinitro-2-methylphenol	43000 UD	43000
(continued)		

Client ID: S8 Lab ID (HSN): 15290 DL Matrix: SOIL Filename: 4067L21

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	17000 UD	17000
4-Bromophenyl-phenylether	17000 UD	17000
Hexachlorobenzene	17000 UD	17000
Pentachlorophenol	43000 UD	43000
Phenanthrene	3200 JD	17000
Anthracene	17000 UD	17000
Carbazole	17000 UD	17000
Di-n-butylphthalate	7000 JDB	17000
Fluoranthene	5400 JD	17000
Pyrene	8700 JD	17000
Butylbenzylphthalate	9100 JD	17000
3,3'-Dichlorobenzidine	17000 UD	17000
Benz(a)anthracene	4000 JD	17000
Chrysene	4600 JD	17000
bis(2-Ethylhexyl)phthalate	9300 JD	17000
Di-n-octylphthalate	.17000 UD	17000
Benzo(b) fluoranthene	7600 JD	17000
Benzo(k) fluoranthene	2500 JD	17000
Benzo(a)pyrene	4900 JD	17000
Indeno(1,2,3-cd)pyrene	7500 JD	17000
Dibenz(a,h)anthracene	17000 UD	17000
Benzo(g,h,i)perylene	7200 JD	17000

Surrogate Recovery		QC LIMITS
2-Fluorophenol	57 % JD	25-121*
Phenol-d5	65 % JD	24-113*
2-Chlorophenol-d4	61 % JD	20-130%
1,2-Dichlorobenzene-d4	54 % JD	20-130%
Nitrobenzene-d5	56 % JD	23-120%
2-Fluorobiphenyl	· 76 % JD	30-115%
2,4,6-Tribromophenol	61 % JD	19-122%
Terphenyl-d14	108 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition.

TT____1 HPN: 2757

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Client ID: S9 Lab ID (HSN): 15291
Matrix: SOIL Filename: 4068K15

Date Sampled: 02/22/94 Sample Size: 15.1 grams
Date Received: 02/24/94 Extract Vol.: 500 uL

Date Extracted: 03/03/94 Dil. Factor: 1
Date Analyzed: 03/10/94 GPC Factor: 2
% Moisture: 14.95

Compounds:	ug/Kg (PPB)	EQL
Phenol	780 U	780
bis(2-Chloroethyl)ether	780 U	780
2-Chlorophenol	780 U	780
1,3-Dichlorobenzene	780 Ŭ	780
1,4-Dichlorobenzene	780 U	780
1,2-Dichlorobenzene	780 U	780
2-Methylphenol	780 U	780
2,2'-oxybis(1-Chloropropane)	780 U	. 780
4-Methylphenol	780 U	780
N-Nitroso-di-n-propylamine	780 U	780
Hexachloroethane	780 U	780
Nitrobenzene	780 U	780
Isophorone	780 U	780
2-Nitrophenol	780 U	780
2,4-Dimethylphenol	780 U	780
bis (2-Chloroethoxy) methane	780 U	780
2,4-Dichlorophenol	780 Ŭ	780
1,2,4-Trichlorobenzene	110 J	780
Naphthalene	780 U	780
4-Chloroaniline	780 U	780
Hexachlorobutadiene	780 U	780
4-Chloro-3-methylphenol	780 U	780
2-Methylnaphthalene	90 J	780
Hexachlorocyclopentadiene	780 U	780
2,4,6-Trichlorophenol	780 U	780
2,4,5-Trichlorophenol	1900 U	1900
2-Chloronaphthalene	780 U	780
2-Nitroaniline	1900 U	1900
Dimethylphthalate	780 U	780
Acenaphthylene	780 U	780
2,6-Dinitrotoluene	780 U	780
3-Nitroaniline	1900 U	1900
Acenaphthene	780 U	780
2,4-Dinitrophenol	1900 U	1900
4-Nitrophenol	1900 U	1900
Dibenzofuran	780 U	780
2,4-Dinitrotoluene	780 U	
Diethylphthalate	780 U	
4-Chlorophenyl-phenylether	780 U	780
Fluorene	110 J	780
4-Nitroaniline	1900 U	. 1900
4,6-Dinitro-2-methylphenol	1900 U	1900

(continued)

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Client ID: S9
Matrix: SOIL

Lab ID (HSN): 15291 Filename: 4068K15

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	780 U	780
4-Bromophenyl-phenylether	780 Ŭ	780
Hexachlorobenzene	780 U	780
Pentachlorophenol	1900 U	1900
Phenanthrene	1100	780
Anthracene	190 J	780
Carbazole	780 Ŭ	780
Di-n-butylphthalate	620 JB	780
Fluoranthene	1500	780
Pyrene	4100 J Y	780
Butylbenzylphthalate	780 U Y	
3,3'-Dichlorobenzidine	780 U Y	
Benz(a)anthracene	1500 T Y	
Chrysene	1500 7 Y	780
bis(2-Ethylhexyl)phthalate	4300 ± Y	780
Di-n-octylphthalate	780 Ŭ	780
Benzo(b) fluoranthene	2800	780
Benzo(k) fluoranthene	860	780
Benzo(a) pyrene	1400	780
Indeno(1,2,3-cd)pyrene	1500	780
Dibenz(a,h)anthracene	780 U	780
Benzo(g,h,i)perylene	1400	780

Surrogate Recovery		QC LIMITS
2-Fluorophenol	30%	25-121%
Phenol-d5	41%	24-113%
2-Chlorophenol-d4	45%	20-130%
1,2-Dichlorobenzene-d4	61%	20-130%
Nitrobenzene-d5	42%	23-120%
2-Fluorobiphenyl	75%	30-115%
2,4,6-Tribromophenol	22*	19-122*
Terphenyl-d14	92%	Y 18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

Note: All results are reported on a 1:y weight basis.

Reference: "EPA Test Methods for Eva. .ating Solid Waste", SW-846,

November 1986, 3rd Edit: 5

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit (estimated value)

^{3 =} Also detected in the associated Blank

Associated internal standard falled method criteria

Client ID: S9 Lab ID (HSN): 15291 DL Matrix: SOIL Filename: 4067L22

Date Sampled: 02/22/94 Sample Size: 15.1 grams
Date Received: 02/24/94 Extract Vol.: 500 uL
Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2

GPC Factor: 2
% Moisture: 14.95

Compounds:	ug/Kg (PPB)	EQL
Phenol	7800 UD	7800
bis(2-Chloroethyl)ether	7800 UD	7800
2-Chlorophenol	7800 UD	7800
1,3-Dichlorobenzene	7800 UD	7800
1,4-Dichlorobenzene	7800 UD	7800
1,2-Dichlorobenzene	7800 UD	7800
2-Methylphenol	7800 UD	7800
2,2'-oxybis(1-Chloropropane)	7800 UD	7800
4-Methylphenol	7800 UD	7800
N-Nitroso-di-n-propylamine	7800 UD	7800
Hexachloroethane	7800 UD	7800
Nitrobenzene	7800 UD	7800
Isophorone	7800 UD	7800
2-Nitrophenol	7800 UD	7800
2,4-Dimethylphenol	7800 UD	7800
bis(2-Chloroethoxy)methane	7800 UD	7800
2,4-Dichlorophenol	7800 UD	7800
1,2,4-Trichlorobenzene	7800 UD	7800
Naphthalene	7800 UD	7800
4-Chloroaniline	7800 UD	7800
Hexachlorobutadiene	7800 UD	7800
4-Chloro-3-methylphenol	7800 UD ·	7800
2-Methylnaphthalene	7800 UD	7800
Hexachlorocyclopentadiene	7800 UD	7800
2,4,6-Trichlorophenol	7800 UD	7800
2,4,5-Trichlorophenol	19000 UD	19000
2-Chloronaphthalene	7800 UD	7800
2-Nitroaniline	19000 UD	19000
Dimethylphthalate	7800 UD	7800
Acenaphthylene	7800 UD	7800
2,6-Dinitrotoluene	7800 UD	7800
3-Nitroaniline	19000 UD	19000
Acenaphth ene	7800 UD	7800
2,4-Dinitrophenol	19000 UD	19000
4-Nitrophenol	19000 UD	19000
Dibenzofuran	7800 UD	7800
2,4-Dinitrotoluene	7800 UD	7800
Diethylphthalate	7800 UD	7800
4-Chlorophenyl-phenylether	7800 UD	7800
Fluorene	7800 UD	7800
4-Nitroaniline	19000 UD '	19000
4,6-Dinitro-2-methylphenol	19000 UD	19000
4,6-Dinitro-2-methylphenol	19000 UD	19000

(continued)

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Client ID: S9 Lab ID (HSN): 15291 DL Matrix: SOIL Filename: 4067L22

	4	
Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	7800 UD	7800
4-Bromophenyl-phenylether	7830 UD	7800
Hexachlorobenzene	7800 UD	7800
Pentachlorophenol	19000 UD	19000
Phenanthrene	1400 JD	7800
Anthracene	7800 UD	7800
Carbazole	7800 UD	7800
Di-n-butylphthalate	1300 JDB	7800
Fluoranthene	2600 JD	7800
Pyrene	5200 JD	7800
Butylbenzylphthalate	7800 UD	7800
3,3'-Dichlorobenzidine	7800 UD	7800
Benz(a)anthracene	1600 JD	. 7800
Chrysene	1900 JD	7800
bis(2-Ethylhexyl)phthalate	9800 D	7800
Di-n-octylphthalate	7800 UD	7800
Benzo(b) fluoranthene	3100 JD	7800
Benzo(k) fluoranthene	860 JD	7800
Benzo(a) pyrene	1400 JD	7800
Indeno(1,2,3-cd)pyrene	1800 JD	7800
Dibenz(a,h)anthracene	7800 UD	7800
Benzo(g,h,i)perylene	1700 JD	7800

Surrogate Recovery 2-Fluorophenol Phenol-d5 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4 Nitrobenzene-d5 2-Fluorobiphenyl 2,4,6-Tribromophenol	39%JD 55%JD 53%JD 58%JD 61%JD 78%JD 24%JD	QC LIMITS 25-121* 24-113* 20-130* 20-130* 23-120* 30-115* 19-122*
Terphenyl-d14	115 % JD	18-137%

TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)

EQL = Estimated Quantitation Limit (lower calibration limit)

U = Undetected at the given EQL

J = Detected below the EQL (estimated value)

E = Exceeds the upper calibration limit estimated value)

B = Also detected in the associated Blank

D = Analysis at a secondary Dilution factor

Note: All results are reported on a ::y weight basis.

Reference: "EPA Test Methods for Eva...: ing Solid Waste", SW-846,

November 1986, 3rd Edit:

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Client ID: S10 Lab ID (HSN): 15292 DL Matrix: SOIL Filename: 4067L23

Date Sampled: 02/22/94 Sample Size: 15.2 grams
Date Received: 02/24/94 Extract Vol.: 2000 uL

Date Extracted: 03/03/94 Dil. Factor: 10
Date Analyzed: 03/09/94 GPC Factor: 2
% Moisture: 40.38

ug/Kg (PPB) Compounds: EQL 44000 UD 44000 Phenol bis(2-Chloroethyl)ether 44000 UD 44000 44000 UD 44000 2-Chlorophenol 44000 UD 1,3-Dichlorobenzene 44000 44000 UD 44000 1,4-Dichlorobenzene 44000 UD 44000 1,2-Dichlorobenzene 2-Methylphenol 44000 UD 44000 44000 UD 44000 2,2'-oxybis(1-Chloropropane) 44000 UD 44000 4-Methylphenol 44000 UD 44000 N-Nitroso-di-n-propylamine 44000 UD 44000 Hexachloroethane 44000 UD 44000 Nitrobenzene 44000 UD 44000 Isophorone 44000 UD 44000 2-Nitrophenol 44000 UD 2,4-Dimethylphenol 44000 44000 UD 44000 bis(2-Chloroethoxy)methane 44000 UD 44000 2,4-Dichlorophenol 44000 UD 44000 1,2,4-Trichlorobenzene 54000 D 44000 Naphthalene 44000 UD 44000 4-Chloroaniline 44000 UD · Hexachlorobutadiene 44000 44000 UD 44000 4-Chloro-3-methylphenol 230000 D 44000 2-Methylnaphthalene 44000 UD 44000 Hexachlorocyclopentadiene 44000 UD 44000 2,4,6-Trichlorophenol 110000 UD 110000 2,4,5-Trichlorophenol 2-Chloronaphthalene 44000 UD 44000 110000 UD 110000 2-Nitroaniline 44000 UD 44000 Dimethylphthalate 44000 UD 44000 Acenaphthylene 44000 UD 44000 2,6-Dinitrotoluene 110000 UD 110000 3-Nitroaniline 14000 JD 44000 Acenaphthene 110000 UD 110000 2,4-Dinitrophenol 110000 UD 110000 4-Nitrophenol 44000 UD 44000 Dibenzofuran 44000 UD 44000 2,4-Dinitrotoluene 44000 UD 44000 Diethylphthalate 44000 UD 44000 4-Chlorophenyl-phenylether 44000 32000 JD Fluorene 110000 UD 110000 4-Nitroaniline 110000 UD 110000 4,6-Dinitro-2-methylphenol

(continued)

Client ID: S10 Lab ID (HSN): 15292 DL Matrix: SOIL Filename: 4067L23

Compounds:	ug/Kg (PPB)	EQL
N-Nitrosodiphenylamine	44000 UD	44000
4-Bromophenyl-phenylether	44000 UD	44000
Hexachlorobenzene	44000 UD	44000
Pentachlorophenol	110000 UD	110000
Phenanthrene	88000 D	44000
Anthracene	7500 JD	44000
Carbazole	44000 UD	44000
Di-n-butylphthalate	44000 UD	44000
Fluoranthene	35000 JD	44000
Pyrene	62000 D	44000
Butylbenzylphthalate	44000 UD	44000
3,3'-Dichlorobenzidine	44000 UD	44000
Benz(a)anthracene	20000 JD	44000
Chrysene	20000 JD	44000
bis(2-Ethylhexyl)phthalate	44000 UD	44000
Di-n-octylphthalate	44000 UD	44000
Benzo(b)fluoranthene	22000 JD	44000
Benzo(k)fluoranthene	7700 JD	44000
Benzo(a)pyrene	15000 JD	44000
Indeno(1,2,3-cd)pyrene	14000 JD	44000
Dibenz(a,h)anthracene	44000 UD	44000
Benzo(g,h,i)perylene	14000 JD	44000

Surrogate Recovery 2-Fluorophenol	40 % JD	QC LIMITS 25-121%
Phenol-d5	69 % JD	24-113%
2-Chlorophenol-d4	61 % JD	20-130%
1,2-Dichlorobenzene-d4	76 % JD	20-130%
Nitrobenzene-d5	156 % JD	23-120%
2-Fluorobiphenyl	92 % JD	30-115%
2,4,6-Tribromophenol	88 % JD	19-122%
Terphenyl-d14	132 % JD	18-137%

- TCL = Target Coumpound List EPA Contract Laboratory Program (OLM01)
- EQL = Estimated Quantitation Limit (lower calibration limit)
 - U = Undetected at the given EQL
 - J = Detected below the EQL (estimated value)
 - = Exceeds the upper calibration limit (estimated value)
 - 3 Also detected in the associated Blank
 - D = Analysis at a secondary Dilution factor

Note: All results are reported on a dry weight basis.

Reference: "EPA Test Methods for Evaluating Solid Waste", SW-846,

November 1986, 3rd Edition

Huntingdon



ecology and environment, inc.

111 WEST JACKSON BLVD CHICAGO, ILLINOIS 60604 TEL 312-663-9415 International Specialists in the Environment

MEMORANDUM

DATE: March 24, 1994

TO: John Nordine, Project Manager, E & E, Chicago, IL

FROM: David Hendren, TAT-Chemist, E & E, Chicago, IL

THRU: Pat Zwilling, ATATL, E & E, Chicago, IL

SUBJ: Polychlorinateddibenso Dioxins and Furans Data Quality

Assurance Review for Scrap Metal site, Chicago, Cook

County, Illinois

REF: Analytical TDD:T05-9402-807 Project TDD:T05-9402-007

Analytical PAN: EIL0831AAA Project PAN: EIL0831SAA

The data quality assurance review of 4 soil samples collected from the site has been completed. Analysis for polychlorinated dibenzodioxins (PCDD) and polychlorinated dibenzofurans (PCDF) was performed by Twin City Testing Corporation (Huntingdon) of St. Paul, Minnesota following U.S. EPA SW-846 Method 8290.

The samples were numbered as follows in the field. The corresponding laboratory identification number is provided:

TAT Sample #	corresponds	to	-	>	Laboratory Sample #
51	•				15270
S2					15277
S3					15278
S8					15290

Data Qualifications:

I. Holding Time: Acceptable

The samples were collected on 2/22/94, extracted on 3/4/94 and analyzed on 3/7/94 and 3/9/94. The holding time criteria of 6 months from sample collection to extraction was satisfied. All

samples were analyzed within the 40 day limit following extraction.

II. Instrument Performance: Acceptable

Analysis of the calibration check solution showed that the % valley between tetrachlorodibenzodioxin (TCDD) isomers (1,2,3,4-) and (2,3,7,8-) was less than 25 %, as required.

III. Calibration:

A. Initial Calibration: Acceptable

The percent relative standard deviations (*RSD) for all analytes in the initial calibration (performed 11/9/93) was less than 15%, as required, except for OCDD (15.66%). Qualification was not judged to be necessary. All specific ion ratio criteria were achieved and signal to noise criteria were satisfied.

B. Continuing Calibration: Acceptable

The continuing calibration standard was analyzed before and following sample analysis. All percent difference of response factors were less than 30%, except for OCDD(34.1%) in the final calibration check run on 3/7/94. Since the initial calibration check was in control and all other parameters were acceptable, qualification was not judged to be necessary.

IV. Method Blank: Acceptable

A method blank was analyzed for each matrix analyzed. HxCDF, HpCDD, and OCDD were detected at 0.39 PPT, 4.4 PPT and 21.0 PPT respectively. All reported values in the samples for these analytes exceeded these levels (over five times) and qualification was not required. No other target analytes or contaminants were detected above the detection limit.

V. Internal Standards: Acceptable

The recoveries of the 13C-internal standards were acceptable in all samples.

VI. Matrix Spike/Matrix Spike Duplicate (MS/MSD): Acceptable

The percent recoveries for the Matrix Spike/Matrix Spike Duplicate (MS/MSD) were within the established quality control limits.

VII. Overall Assessment of Data For Use:

The overall usefulness of the data is based on the criteria outlined in "Quality Assurance/Quality Control Guidance for Removal Activities" (OSWER 9360.4-01 April, 1990). Based upon the information provided, the data are acceptable for use.

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*TWIN CITY TESTING CORPORATION*
*METHOD 8290 ANALYSIS RESULTS *
Client....ECOLOGY & ENVIRONMENT
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Client's Sample ID.....S1 (1:2.5 DILUTION)
TCT Sample ID.......15270
Analysis Date......3/7/94 12:35
Filename.......V40307D
Injected By.......MJK
Total Amount Extracted...0.0129 kg CCAL Filename......V40307B
Method Blank ID.....BLANK-15828 Extraction Date.....3/4/94

NATIVE ISOMERS	CONC. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF TOTAL TCDF	220 * 1500		2378-TCDF-13C 2378-TCDD-13C 12378-PeCDF-13C	2.00 2.00 2.00	105 98 87
2378-TCDD TOTAL TCDD	ND 34	11	23478-PeCDF-13C 12378-PeCDD-13C 123478-HxCDF-13C.	2.00 2.00 2.00	91 81 102
12378-PeCDF 23478-PeCDF TOTAL PeCDF	94 260 1300		123678-HXCDF-13C. 234678-HXCDF-13C. 123789-HXCDF-13C. 123478-HXCDD-13C.	2.00 2.00 2.00 2.00	81 93 81 98
12378-PeCDD TOTAL PeCDD	ND ND	310	123678-HxCDD-13C. 1234678-HpCDF-13C 1234789-HpCDF-13C	2.00 2.00 2.00	97 73 72
123478-HxCDF 123678-HxCDF 234678-HxCDF 123789-HxCDF	250 160 230 95		1234678-HPCDD-13C OCDD-13C	2.00	82 82 NA
TOTAL HXCDF	2000		123789-HxCDD-13C.	2.00	NA
123478-HxCDD 123678-HxCDD 123789-HxCDD	21 70 33		2378-TCDD-37C14	0.20	INT
TOTAL HXCDD	240		Total 2378-TCDD Equivalence:	265	ng/kg
1234678-HpCDF 1234789-HpCDF TOTAL HpCDF	900 120 1500		(Using ITE Fact	ors/DB-5	Dátas
1234678-HpCDD TOTAL HpCDD	730 1400				
OCDF OCDD	770 4300				

^{*} Value may include contributions from other TCDF isomers.

All values are expressed on a dry weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.)
LOD = Limit of Detection
ND = Not Detected
NA = Not Applicable
INT = Interference

INT = Interference

TCT Invoice Number....4411 94-2757

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*METHOD 8290 ANALYSIS RESULTS *
                              Client.... ECOLOGY & ENVIRONMENT
Client's Sample ID.....S2 (1:20 DILUTION)
TCT Sample ID......15277
Analysis Date......3/9/94 14:42
Extraction Date.....3/4/94
NATIVE
                     CONC.
                                                        INTERNAL
                                                                           ng's
                                                                                     PERCENT
ISOMERS
                                                        STANDARDS
                                                                          ADDED
                     ng/kg
                                   ng/kg
                                                                                     RECOVERY
2378-TCDF
                                                 2378-TCDF-13C....
                                   10000
                                                                             2.00
                                                 2378-TCDF-13C....

2378-TCDD-13C....

12378-PeCDF-13C...

23478-PeCDD-13C...

12378-PeCDD-13C...

123478-HxCDF-13C...

234678-HxCDF-13C...

234678-HxCDF-13C...

123789-HxCDF-13C...
                                                                                          103
TOTAL TCDF
                     17000
                                                                             2.00
                                                                                            60
                                                                             2.00
                                                                                            89
2378-TCDD
                         ND
                                       22
                                                                                          120
TOTAL TCDD
                        580
                                                                             2.00
                                                                                            97
                                                                             2.00
                                                                                           128
                                                                             2.00
12378-PeCDF
                      4400
                                                                                            88
23478-PeCDF
                                                                             2.00
                       6300
                                                                                            94
TOTAL PeCDF
                                                                             2.00
                     42000
                                                                                            98
                                                                             2.00
                                                                                            98
                                                 123678-HxCDD-13C.
1234678-HpCDF-13C
1234789-HpCDF-13C
                                   4200
12378-PeCDD
                         ND
                                                                             2.00
                                                                                            90
TOTAL PeCDD
                         ND
                                   ----
                                                                             2.00
                                                                                            95
                                                                             2.00
                                                                                            71
123478-HxCDF
                         ND
                                   20000
                                                 1234678-HPCDD-13C
                                                                             2.00
                                                                                            72
123678-HxCDF
                                                 OCDD-13C.....
                                    4100
                         ND
                                                                             4.00
234678-HXCDF
123789-HXCDF
                      2700
                                                 1234-TCDD-13C....
123789-HxCDD-13C.
                                                                             2.00
                        890
                                                                                           NA
TOTAL HXCDF
                     15000
                                                                             2.00
                                                                                           NA
123478-HxCDD
123678-HxCDD
123789-HxCDD
                         ND
                                     140
                                                 2378-TCDD-37C14..
                                                                             0.20
                                                                                          INT
                        250
130
TOTAL HXCDD
                        380
                                                     Total 2378-TCDD
                                                                             4004 ng/kg
                                                     Equivalence:
                     12000
                                                     (Using ITE Factors/DB-5 Data)
1234678-HpCDF
1234789-HDCDF
TOTAL HDCDF
                      3600
                                   ____
                     16000
1234678-HpCDD
TOTAL HpCDD
                      2900
                                   ____
                      5900
OCDF
                     30000
                     22000
All values are expressed on a dry weight basis.
CONC= Concentration (Totals include 2378-substituted isomers.)
LOD = Limit of Detection
ND = Not Detected
NA = Not Applicable
INT = Interference
                                          TCT Invoice Number....4411 94-2757
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TWIN CITY TESTING CORPORATION

Huntingdon

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*TWIN CITY TESTING CORPORATION*
                          *METHOD 8290 ANALYSIS RESULTS *
                           Client....ECOLOGY & ENVIRONMENT
Client's Sample ID.....S3 (1:2.5 DILUTION) TCT Sample ID.........15278
Analysis Date.....3/7/94 14:30 Filename.....V40307F Injected By......MJK
Total Amount Extracted...0.0155 kg
Extraction Date......3/4/94
                                                                     ng's
                                                   INTERNAL
                                                                              PERCENT
                                LOD
NATIVE
                   CONC.
                                                                    ADDED
                                                   STANDARDS
                                                                              RECOVERY
ISOMERS
                   ng/kg
                               ng/kg
                                            2378-TCDF-13C...

2378-TCDD-13C...

12378-PeCDF-13C..

23478-PeCDF-13C..

12378-PECDD-13C..

123478-HxCDF-13C.

123678-HxCDF-13C.

234678-HxCDF-13C.
                                                                       2.00
2378-TCDF
                     1700 *
TOTAL TCDF
                                                                       2.00
                                                                                   111
                    9400
                                                                       2.00
                                                                                   116
2378-TCDD
TOTAL TCDD
                       ND
                                                                      2.00
                                                                                   124
                                   18
                                                                       2.00
                      130
                                                                                   106
                                                                      2.00
                                                                                   120
12378-PeCDF
                     550
                                                                       2.00
                                                                                    91
                                                                                   114
95
23478-PeCDF
                     1300
                                                                       2.00
                                             123789-HxCDF-13C.
                                                                      2.00
TOTAL PeCDF
                     7600
                                             123478-HxCDD-13C.
                                                                      2.00
                                                                                   114
                                             123678-HXCDD-13C.
1234678-HpCDF-13C
12378-PeCDD
                       ND
                                  160
                                                                       2.00
                                                                                    99
                                                                                    87
TOTAL PeCDD
                                                                       2.00
                       ND
                                             1234789-HPCDF-13C
                                                                       2.00
                                                                                    84
                                                                      2.00
                                                                                    87
                                             1234678-HpCDD-13C
123478-HxCDF
                      810
                                             ocdd-13c.....
                                                                      4.00
                                                                                    78
123678-HxCDF
                      640
                      990
220
234678-HxCDF
123789-HxCDF
                                             1234-TCDD-13C...
                                                                      2.00
                                                                                    NA
                                                                      2.00
                                             123789-HxCDD-13C.
                                                                                    NA
TOTAL HXCDF
                    7700
                      100
170
                                                                                   INT
                                             2378-TCDD-37C14..
                                                                      0.20
123478-HxCDD
123678-HxCDD
123789-HxCDD
                      120
                                                Total 2378-TCDD
TOTAL HXCDD
                    1100
                                                Equivalence:
                                                                      1207
                                                                             ng/kg
                                                 (Using ITE Factors/DB-5 Data)
                    3100
1234678-HpCDF
1234789-HDCDF
                      330
                     4500
TOTAL HPCDF
1234678-HpCDD
TOTAL HpCDD
                    1500
                    2900
OCDF
                     1600
OCDD
                    3900
* Value may include contributions from other TCDF isomers.
All values are expressed on a dry weight basis.
CONC= Concentration (Totals include 2378-substituted isomers.)
LOD = Limit of Detection
ND = Not Detected
    = Not Applicable
                                      TCT Invoice Number....4411 94-2757
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Huntingdon

INT = Interference

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*TWIN CITY TESTING CORPORATION*
                            *METHOD 8290 ANALYSIS RESULTS *
                             Client....ECOLOGY & ENVIRONMENT
Client's Sample ID.....S8 (1:20 DILUTION) TCT Sample ID......15290
Analysis Date.....3/9/94 13:41 Filename.....V40309F Injected By......MJK
Total Amount Extracted...0.0144 kg
CCAL Filename......V40309B
Method Blank ID.....BLANK-15828
Extraction Date.....3/4/94
                                                                          ng's
NATIVE
                     CONC.
                                   LOD
                                                       INTERNAL
                                                                                   PERCENT
ISOMERS
                                                       STANDARDS
                                                                         ADDED
                                                                                   RECOVERY
                     ng/kg
                                  ng/kg
                                                2378-TCDF-13C....
2378-TCDD-13C....
12378-PeCDF-13C...
2378-TCDF
TOTAL TCDF
                                                                                          87
                         ND
                                    380
                      1900
                                                                            2.00
                                                                                          79
                                                                            2.00
                                                                                          99
                                                123/8-PECDF-13C..
23478-PECDF-13C..
12378-PECDD-13C..
123478-HxCDF-13C.
                                                                                        132
93
105
75
                                                                           2.00
2378-TCDD
TOTAL TCDD
                        ND
                                      17
                       150
                                                                            2.00
                                                123678-HXCDF-13C.
                        ND
                                  6000
                                                                           2.00
12378-PeCDF
                                                234678-HXCDF-13C.
123789-HXCDF-13C.
123478-HXCDD-13C.
                                                                                          91
87
23478-PeCDF
TOTAL PeCDF
                       220
                                                                           2.00
                      3300
                                                                                          92
                                                123678-HxCDD-13C.
1234678-HpCDF-13C
                                                                            2.00
                                                                                          86
12378-PeCDD
TOTAL PeCDD
                         ND
                                      31
                                                                                          83
67
                                                                            2.00
                        ND
                                                1234789-HPCDF-13C
1234678-HPCDD-13C
                                                                            2.00
                                                                                          77
                        500
                                                                            2.00
123478-HxCDF
                                    340
                                                OCDD-13C....
123678-HxCDF
                        ND
                                                                           4.00
                                                                                          60
234678-HXCDF
123789-HXCDF
                       360
170
                                    ___
                                                1234-TCDD-13C....
123789-HxCDD-13C.
                                                                           2.00
                                                                                          NA
                      2700
                                                                           2.00
TOTAL HXCDF
                                                                                          NA
                                                                           0.20
123478-HxCDD
                         29
                                                2378-TCDD-37Cl4..
                                                                                        INT
123678-HxCDD
                       670
                                  ----
123789-HxCDD
                       190
TOTAL HXCDD
                      1800
                                                    Total 2378-TCDD
                                                    Equivalence:
                                                                                   ng/ka
                                                                             360
                                                    (Using ITE Factors/DB-5 Data)
                      1100
1234678-HpCDF
1234789-HpCDF
                       240
TOTAL HPCDF
                      1300
                      2800
1234678-HpCDD
TOTAL HPCDD
                      5700
                                  ____
OCDF
                      1600
                     15000
OCDD
                                  ____
All values are expressed on a dry weight basis.
CONC= Concentration (Totals include 2378-substituted isomers.)
LOD = Limit of Detection
ND = Not Detected
    = Not Applicable
NA
                                         TCT Invoice Number....4411 94-2757
INT = Interference
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APPENDIX C RCMS COST ESTIMATE

Summery Report

Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: ILO831SA Date: 05/02/94

Cleanup Contractor: ESE5 - ESE TAT Contractor: ECOLOGY&ENVIRON

Cost Projection Summery ****************

Contractor Personnel	282,065,90
Contractor Equipment	55,316.98
Unit Rate Materials	115,944.00
At Cost Materials	1,302.50
Subcontractors	14,220.00
Waste Transportation	7,312.50
Waste Disposal	702,000.00
Cleanup Contractor Subtotal	1,178,161.88
Federal and State Agencies	0.00
Extramural Subtotal	1,178,161.88
15 % Extramural Contingency	176,724.28
Extramural Subtotal	1,354,886.16
TAT Personnel	97,632.00
TAT Special Projects	0.00
TAT Analytical Services	0.00
Total TAT Costs	97,632.00
Other Cost Items	0.00
and the second	
Extramural Subtotal	1,452,518.16
10 % Project Contingency	145,251.82
Total Extramural Cost	1,597,769.98
EPA Regional Personnel	32,400.00
EPA Non-Regional Personnel	0.00
EPA Headquarters Direct	0.00
(0 % of Regional Hours)	
EPA Indirect	65,880.00
EPA Total	98,280.00
Acciona Food	
Project Total	1,696,049.98

Summery Report (cont.)

Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: IL0831SA

Cleanup Contractor: ESE5 - ESE

Date: 05/02/94

TAT Contractor: ECOLOGY&ENVIRON

Project Scope

		Estimated	
Number	Step/Milestone	Duration	Cost
•••••		• • • • • • • • • • • • • • • • • • • •	
000	GENERAL SITE COSTS	120 Days	1,696,049.98
999	ARCHIVE COSTS	120 Days	0.00
	•		1.696.049.98

9,047.70

Detailed Report By Category Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: IL0831SA

Cleanup Contractor: ESE5 - ESE TAT Contractor: ECOLOGY&ENVIRON

Cost Projection Detail - By Category 118488888881118888888888 1184828888

Cont	ract	100	Pers	ionnel

Job Category	Number of Employees	Number of Days	Hrs per Day	Labor	PD, Lodge Travel	Total Charge
***************************************					••••••	•••••
000 - GENERAL SITE COSTS						
						.=
S105-Response Manager	1	90 ·	12.00	52,023.60	13,355.10	65,378.70
S205-Equipment Operato	1	90	12.00	41,047.20	13,355.10	54,402.30
S210-field Clerk/Typis	1	90	12.00	28,087.20	13,355.10	41,442.30
S215-Laborer	2	90	12.00	61,934.40	26,710.20	88,644.60
S620-Level C	4	50	12.00	2,520.00	29,678.00	32,198.00
						••••••
			Total for	GENERAL SITE	COSTS :	282,065.90

Contractor Equipment

urractor Editibular								
Equipment Name	Number Needed	Reg Days	Hours /day	Stby Days	Hob/Demob Days	Decon Days	Hileage	Total Charge
000 - GENERAL SITE COSTS	•••••	••••	•••••	•••••	*********	•••••	• • • • • • • • • • • • • • • • • • • •	• • • • • • • • • • • • • • • • • • • •
10320-Box-2 ton	1	90	12.00	0	4	0	H/A	4,075.36
13610-Pickup-2 wheel drive	1	90	12.00	0	4	0	N/A	2,695.05
15430-Van-Passenger	1	90	12.00	0	4	0	N/A	4,136.93
22020-Decon-8x25	1	90	12.00	0	4	0	N/A	1,948.70
25520-Office-8x20	1	90	12.00	0	4	0	N/A	1,145.27
33055-Excavator-CAT 215C	1	90	12.00	. 0	4	0	N/A	21,208.42
35530-Loader/Track-CAT 943	1	90	12.00	0	4	0	N/A	19,068.20
72320-Computer-Portable PC	1	90	12.00	0	4	0	N/A	1,039.05

Total for GENERAL SITE COSTS : 55,316.98

Total equipment cost: 55,316.98

Total personnel cost: 282,065.90

Date: 05/02/94

Unit Rate Materials

	Material Name	Material Use	Unit Cost	Number of Units	Total Charge
000	GENERAL SITE COSTS				
	TOPSOIL	COVER/FILL	4.000	4550.0 CU/YD	18,928.00
	PCB TEST KITS	SAMPLING	120.000	25.0 KIT	3,000.00
	PPE	H & S	75.000	1200.0 EACH	93,600.00

Page:

Detailed Report By Category (cont.) Initial Cost Projection Scenario: SCRAP METAL

Date: 05/02/94

7,312.50

3.25 10

Cleanup Contractor: ESE5 - ESE TAT Contractor: ECOLOGY&ENVIRON

Projection ID Number: [L0831SA

PCB METAL SOILS

Unit Rate Materials Material Name Material Use Unit Cost Number of Units ______ 2.000 200.0 JARS SAMPLE JARS SAMPL ING 416.00 Total for GENERAL SITE COSTS : 115,944,00 Total unit rate materials cost: At Cost Materials Total Material Use Quantity/Amount Material Name Charge 000 - GENERAL SITE COSTS 500 FUEL DEISEL 1.25 500 FUEL GASOLINE 1.25 FIELD SUPPLIES MULTI USE BULK 1,040.00 OFFICE SUPPLIES ADMINISTRATIVE BULK 260.00 Total for GENERAL SITE COSTS : 1,302.50 Total at cost materials cost: 1,302.50 Subcontractors Service Subcontractor Billing 000 - GENERAL SITE COSTS 3.0 MONTHS 1,200.00 ELECTRIC SERVICE FD I SOM 832.00 4.0 MONTH ILLINOIS BELL PHONE SERVICE 3.0 MONTHS 540.00 PORT-O-POTTY PORT BATHROOM REN 7,072.00 SECURITY SITE WATCHMAN 12.0 WEEKS ANALYTICAL ANALYTICAL DATA 3.0 MONTH 4,160.00 CULICAN WATER DRINKING WATER 3.0 MONTH Total for GENERAL SITE COSTS : 14,220.00 14,220.00 Total subcontractor cost: Waste Transportation Total Cost Waste Type Amount , Loads Per Mile Hiles 000 - GENERAL SITE COSTS

4500 CU/YD 225

Page:

Detailed Report By Category (cont.) Initial Cost Projection Scenario: SCRAP METAL

Projection ID Number: [L0831SA Date: 05/02/94

TAT Contractor: ECOLOGY&ENVIRON Cleanup Contractor: ESE5 - ESE

		_		Cost		Total
	e Type		Loads		Miles	Charge
			Tota	nt for GENERAL	L SITE COSTS :	7,312.5
				Total transp	portation cost:	7,312.5
ste Disposal				Ho of	Unit	Tagal
	•	Disposal Method		Units	Cost	Charge
000 - GENER	AL SITE COST		• • • • • • • • • • • • • • • • • • • •		• • • • • • • • • • • • • • • • • • • •	**********
	TAL SOILS		CU/YD	4500	150.00	-
			Tota	al for GENERA	L SITE COSTS :	702,000.0
					•	•••••
				Total	disposal cost:	702,000.0
deral and Stat	e Agencies			Total	disposal cost:	
deral and Stat	e Agencies				disposal cost:	0.0
deral and Stat T Personnel	-	W W			al Contingency:	0.0 176,724.2
T Personnel	Number of Days		urly ate	15 % Extramur: Labor	al Contingency: PO, Lodge Travel	0.0 176,724.2 Total Charge
T Personnel Level	Number of Days LAL SITE COST	Day R	urly ate	15 % Extramur: Labor	al Contingency: PO, Lodge Travel	0.0 176,724.2 Total Charge
T Personnel Level 000 - GENER	Number of Days LAL SITE COST	Day R	urty ate	15 % Extr am ura	al Contingency: PÔ, Lodge Travel	0.0 176,724.2 Total Charge
T Personnel Level 000 - GENER	Number of Days LAL SITE COST	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	al Contingency: PÔ, Lodge Travel	0.0 176,724.2 Total Charge 97,632.0
T Personnel Level 000 - GENER	Number of Days LAL SITE COST	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	PO, Lodge Travel 0.00 L SITE COSTS :	0.0 176,724.2 Total Charge 97,632.0
T Personnel Level 000 - GENER	Number of Days LAL SITE COST	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	PO, Lodge Travel	0.0 176,724.2 Total Charge 97,632.0
T Personnel Level 000 - GENER	Number of Days AL SITE COST 90	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	PO, Lodge Travel 0.00 L SITE COSTS :	0.0 176,724.2 Total Charge 97,632.0 97,632.0
T Personnel Level 000 - GENER	Number of Days AL SITE COST 90	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	PO, Lodge Travel 0.00 L SITE COSTS :	97,632.0 97,632.0
T Personnel Level 000 - GENER PL3	Number of Days AL SITE COST 90	Day R	urly ate 0.00 9	15 % Extramura Labor 7,632.00	PO, Lodge Travel 0.00 L SITE COSTS :	0.0 176,724.2 Total Charge 97,632.0

Page:

Projection ID Number: IL0831SA

Date: 05/02/94

TAT Contractor: ECOLOGY&ENVIRON Cleanup Contractor: ESE5 - ESE

EPA Regi	onal Personnel Title	Number of Days	•	Hourly Rate	Labor	PD, Lodge Travel	
000	- GENERAL SITE CO	STS					
	osc	90	12.00	30.00	32,400.00	0.00	32,400.00
				Total	for GENERAL SI	TE COSTS :	32,400.00
				•		•••	
				Total EPA	Regional Perso	onnel Cost;	32,400.00
				EPA Headq	warters Cost:		0.00
				(0 % of	Regional hours)	
				EPA Indir	ect Cost:		65,880.00
				(1080 hou	rs 8 \$61.00 pe	r hour)	
EPA Non-	Regional Personnel	l				•	0.00
			T	otal EPA Cos	it:		32,400.00

Total site cost: 1,630,169.98



MEMORANDUM

November 9, 1994 DATE:

Raghu Nagam, Project Manager, E & E, Chicago, IL TO:

Herbert B. Langer, TAT-Chemical Engineer, E & E, NC FROM:

Detroit, MI

THRU: Dave Hendren, TAT-Chemist, E & E, Chicago, IL

Asbestos Analysis Data Quality Assurance Review, Scrap SUBJ:

Metal, Chicago, Cook County, IL

REF: Analytical TDD: T05-9410-808 Project TDD: T05-9410-143

Analytical PAN: EIL0831ABA Project PAN: EIL0831FAA

The data quality assurance review for the fibrous sample collected from the Scrap Metal site in Chicago, Illinois, has been completed. Determination of Asbestos in bulk samples by Polarized Light Microscopy was performed by Huntington Environmental Inc., St. Paul, Minnesota. sample was labeled S-17, corresponding to laboratory identification number 43017.

Data Qualifications

Holding Time: Acceptable Ι

> The sample was collected October 17, 1994, and analyzed upon receipt by the laboratory. The OSWER Directive 9360.4-01 does not include criteria regarding holding times for this method.

ΙI Overall Assessment of Data for Use: Acceptable

There are no criteria specified in the U.S. EPA OSWER Directive 9360.4-01 for the determination of asbestos in samples. The analytical method adopted by the laboratory uses visual comparison of the sample material with standard samples to determine the presence, type, and amount of Asbestos in the sample. Based on the analytical method, the data is considered acceptable for use as reported.

PLAINTIFF'S

ANALYTICAL RESULTS

Client ID:

S-17, Trailer

on West Plot

TCT ID:

43017

Sample Identification Asbestos Type & Approximate Percent

Balance of Material & Approximate Percent

Grey fibrous

ND

Glass

90

Non-Fibrous

10

ND = Not Detected

Reference:

The Interim Method of the Determination of Asbestos in Bulk Insulation Samples, EPA

600/M4-82-020, December 1982.



ecology and environment, inc.

12251 UNIVERSAL, TAYLOR, MICHIGAN 48180, TEL. (313) 946-0900 International Specialists in the Environment

MEMORANDUM

November 9, 1994 DATE:

Raghu Nagam, Project Manager, E & E, Chicago, IL TO:

Herbert B. Langer, TAT-Chemical Engineer, E & E, Detroit, MI FROM:

Detroit, MI

Dave Hendren, TAT-Chemist, E & E, Chicago, IL THRU:

Inorganic Data Quality Assurance Review, Scrap Metal, SUBJ:

Chicago, Cook County, IL

Analytical TDD: T05-9410-808 Project TDD: T05-9410-143 REF:

Analytical PAN: EIL0831ABA Project PAN: EIL0831FAA

The data quality assurance review for the six soil samples collected from the Scrap Metal site in Chicago, Illinois, has been completed. Sample analyses were for total Resource Conservation and Recovery Act (RCRA) listed metals following EPA methods 6010 and 7470/71 for mercury. The analyses were performed on sample extracts and Toxicity Characteristic Leaching Procedure (TCLP) leachates. The analyses were performed by Huntington Environmental Inc., St. Paul, The samples were labeled consecutively from S-11 through S-16, corresponding to laboratory identification numbers 43011 through 43016, respectively.

Data Oualifications

Ι Sample Holding Time: Acceptable

> The samples were collected October 17, 1994. Total metal analyses were performed October 21, 1994. Analyses of the TCLP leachates were performed October 24, 1994. All analyses for mercury were performed October 24, 1994. All of these analyses were performed within the required holding times for the methods and matrix.

Initial and Continuing Calibration Verification: Acceptable ΙI

Initial calibration standards and blanks were analyzed at the beginning of each analytical run and after a maximum of ten samples were analyzed, as required. The reported values for the calibration standards were within the required ranges.

III Blanks: Acceptable

A method blank was prepared and analyzed with the samples. None of the target analytes were detected in the blank.

IV Interference Check Sample (ICS) Analysis: Acceptable

A total of two ICSs were run within eight hours of the metals analysis. The ICS results were within \pm 20% of the mean value as required.

V Matrix Spike/Matrix Spike Duplicate (MS/MSD): Acceptable

A MS and MSD were prepared using sample S-12 and the TCLP leachate of sample S-12. The percent recoveries of the spike metals and relative percent differences between MS and MSD results for the actual sample were poor. The laboratory reported that this was due to the debris found in the sample and the difficulties it presented in homogenizing the MS and MSD. The percent recoveries of the spike compounds from the leachate and the relative percent difference between leachate MS and MSD results were acceptable.

VI Overall Assessment of Date for Use

The overall usefulness of the data is based on the criteria outlined in OSWER Directive 9360.4-01 (April 1990), Data Validation Procedures, Section 3.0, Metallic Inorganic Parameters. Based upon the information provided, the data are considered acceptable for use as reported.

METAL RESULTS

(All values are in mg/Kg which is equal to parts-per-million)

Client ID:

S-11, Soil Pile East of Gate of EW Plot S-12, Motor Pile South of S-11 EW Plot

TCT ID:

43011

43012

101 10.	73011	10015			
<u>Parameter</u>			<u>POL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	· <u></u>	25	10/21/94	6010
Arsenic		ND	5.0	10/21/94	6010
Barium	350	330	0.50	10/21/94	6010
Cadmium	44	56	25	10/21/94	6010
Chromium	150	140	25	10/21/94	6010
Lead	4,500	5,100	125	10/21/94	6010
Mercury	25	27	2.0	10/24/94	7471
Selenium	ND	••	25	10/21/94	6010
Selenium		ND	5.0	10/21/94	6010
Silver	48	14	0.50	10/21/94	6010

All results are reported on a dry weight basis.

ND = Not Detected

PQL = Practical Quantitation Limit

Reference:

METAL RESULTS

(All values are in mg/Kg which is equal to parts-per-million)

Client ID:

S-13, Large Ferrous Metal Pile EW Plot S-14, Copper Fines Pile, East of S-13 - EW Plot

Test Test Date <u>Method</u>
/21/94 6010
/21/94 6010
/21/94 6010
/21/94 6010
/21/94 6010
/21/94 6010
/21/94 6010
/24/94 7471
/21/94 6010

4.5

0.50

10/21/94

6010

All results are reported on a dry weight basis.

6.0

ND = Not Detected

Silver

PQL = Practical Quantitation Limit

Reference:

METAL RESULTS

(All values are in mg/Kg which is equal to parts-per-million)

Client ID:

S-15, Sorted Steel Pile, East of S-14- SW Plot S-16, Process Area By Conveyor East Plot

43015 43016 TCT ID: Test Test **PQL** Date Method **Parameter** ND 5.0 Arsenic 10/21/94 6010 14 Barium 83 170 0.50 10/21/94 6010 Cadmium ND 61 25 10/21/94 6010 Chromium 380 320 25 10/21/94 6010 Lead 77 2.5 10/21/94 6010 125 Lead 790 10/21/94 6010 Mercury 27 30 2.0 10/24/94 7471

ND

5.0

All results are reported on a dry weight basis.

ND

5.2

ND = Not Detected

PQL = Practical Quantitation Limit

Reference:

Selenium

Silver

EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

5.0

0.50

10/21/94

10/21/94

6010

6010

TCLP METAL RESULTS

(All values are in μ g/L which is equivalent to parts-per-billion)

Client ID:

Method Blank

LCS

TCT ID:

<u>Parameter</u>			POL	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	87%	100	10/24/94	6010
Barium	ND	91%	10	10/24/94	6010
Cadmium	ND	97%	10	10/24/94	6010
Chromium	ND	97%	10	10/24/94	6010
Lead	ND	99%	50	10/24/94	6010
Mercury	ND	100%	0.40	10/24/94	7470
Selenium	ND	89%	100	10/24/94	6010
Silver	ND	83%	10	10/24/94	6010
TCLP Date:	10/20/94	10/20/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

TCLP METAL RESULTS

(All values are in μ g/L which is equivalent to parts-per-billion)

S-11, Soil Pile

Client ID: East of Gate of EW Plot

OF EW PIO

TCT ID:	43011	MS(43011)	MSD(43011)			
<u>Parameter</u>				<u>POL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	89%	91%	100	10/24/94	6010
Barium	4,100	120%	130%	10	10/24/94	6010
Cadmium	780	91%	93%	10	10/24/94	6010
Chromium	46	90%	89%	10	10/24/94	6010
Lead	37,000	NA^1	NA ¹	50	10/24/94	6010
Mercury	0.50	79%	89%	0.40	10/24/94	7470
Selenium	ND	98%	98%	100	10/24/94	6010
Silver	ND	83%	83%	10	10/24/94	6010
TCLP Date:	10/20/94	10/20/94	10/20/94			

¹ Recovery could not be determined due to the large amount of lead present in the sample matrix.

PQL = Practical Quantitation Limit

ND = Not Detected

NA = Not Applicable

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

TCLP METAL RESULTS

(All values are in μ g/L which is equivalent to parts-per-billion)

Client ID:

S-12, Motor Pile South of S-11 EW Plot

TCT ID:

43012

Parameter		<u>POL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	100	10/24/94	6010
Barium	1,900	10	10/24/94	6010
Cadmium	490	10	10/24/94	6010
Chromium	15	10	10/24/94	6010
Lead	8,500	50	10/24/94	6010
Mercury	ND	0.40	10/24/94	7470
Selenium	ND	100	10/24/94	6010
Silver	ND	10	10/24/94	6010
TCLP Date:	10/20/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

TCLP METAL RESULTS

(All values are in μ g/L which is equivalent to parts-per-billion)

Client ID:	S-13, Large Ferrous Metal Pile EW Plot	S-14, Copper Fines Pile, East of S-13 - EW Plot

TCT ID:	43013	43014			
<u>Parameter</u>			<u>POL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	100	10/24/94	6010
Barium	1,700	1,800	10	10/24/94	6010
Cadmium	ND	84	30	10/24/94	6010
Chromium	ND	ND	10	10/24/94	6010
Lead	72	ND	50	10/24/94	6010
Mercury	ND	ND	0.40	40/24/94	7470
Selenium	ND	ND	100	10/24/94	6010
Silver	ND	ND	10	10/24/94	6010
TCLP Date:	10/20/94	10/20/94			

PQL = Practical Quantitation Limit

ND = Not Detected

Reference: Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.

TCLP METAL RESULTS

(All values are in $\mu g/L$ which is equivalent to parts-per-billion)

Client ID:

S-15, Sorted Steel Pile, East of S-16, Process Area By Conveyor

S-14- SW Plot

East Plot

TCT ID:

43015

43016

101 121					
<u>Parameter</u>			<u>PQL</u>	Test <u>Date</u>	Test <u>Method</u>
Arsenic	ND	ND	100	10/24/94	6010
Barium	1,400	1,800	10	10/24/94	6010
Cadmium	ND	600	30	10/24/94	6010
Chromium	ND	ND	10	10/24/94	6010
Lead	89	63	50	10/24/94	6010
Mercury	ND	ND	0.40	10/24/94	7470
Selenium	ND	ND	100	10/24/94	6010
Silver	ND	ND	10	10/24/94	6010
TCLP Date:	10/20/94	10/20/94			
					

PQL = Practical Quantitation Limit

ND = Not Detected

Reference:

Federal Register, Volume 55, Number 126, June 1990, 40CFR, Method 1311.



MEMORANDUM

DATE: November 9, 1994

TO: Raghu Nagam, Project Manager, E & E, Chicago, IL

FROM: Herbert B. Langer, TAT-Chemical Engineer, E & E, 2

Detroit, MI

THRU: Dave Hendren, TAT-Chemist, E & E, Chicago, IL

SUBJ: Polychlorinated Biphenyl Compound Data Quality

Assurance Review, Scrap Metal, Chicago, Cook County, IL

REF: Analytical TDD: T05-9410-808 Project TDD: T05-9410-143

Analytical PAN: EIL0831ABA Project PAN: EIL0831FAA

The data quality assurance review for the six soil samples collected from the Scrap Metal site in Chicago, Illinois, has been completed. Sample analyses were for polychlorinated biphenyl compounds (PCB) following EPA method 8080. The analyses were performed by Huntington Environmental Inc., St. Paul, Minnesota. The samples were labeled consecutively from S-11 through S-16, corresponding to laboratory identification numbers 43011 through 43016, respectively.

Data Oualifications

I Sample Holding Time: Acceptable

The samples were collected October 17, 1994, extracted October 19, 1994, and analyzed October 23, 1994. All activities were performed within the required holding times for the method and matrix.

II Pesticide Instrument Performance: Acceptable

Standard chromatograms show adequate peak resolution. Surrogate compounds were diluted out of the samples.

III Initial Calibration and Continuing Calibration: Acceptable

Initial calibrations were performed October 22, 1994, using all target Aroclors at five different concentrations. The

percent relative standard deviations between Aroclor response factors were less than twenty, as required by EPA method 8000. Continuing calibrations were performed during the run. Percent differences between initial and continuing calibration response factors were less than fifteen for both the quantitation and confirmation columns, as required.

IV Method Blank: Acceptable

A method blank was prepared and analyzed with the samples. None of the target compounds were detected in the blank.

- V Optional Quality Control Analyses:
 - A. Matrix Spike/Matrix Spike Duplicates (MS/MSD): No Action

A MS and MSD were prepared using the sample S-14. The PCB concentrations in the extract and dilutions required to bring the extract into the instrument's calibration range made spike recoveries impossible.

B. Surrogate Recoveries: No Action

A total of two surrogate compounds were added to the samples, blank, MS, and MSD. The surrogate compounds were diluted out of the samples, MS, and MSD. The surrogate recoveries for the blank were within the laboratory's quality control guidelines

VI Compound Identification: Acceptable

Positive results were identified using correct retention time windows, peak height ratios, and fingerprint patterns. Dual column confirmations of the positive results were performed.

VII Compound Quantitation and Reported Detection Limits: Acceptable

Compound quantitation and detection limits have been correctly adjusted to reflect matrix effects, dry weight factors, and dilutions.

VIII Overall Assessment of Data for Use:

The overall usefulness of the data is based on the criteria outlined in OSWER Directive 9360.4-01 (April 1990), Data Validation Procedures, Section 7.0, PCBs. Based upon the information provided, the data are considered acceptable for use as reported.

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-11, Soil Pile East of Gate of EW Plot

TCT ID:	43011	
Parameter:		POL
PCB 1016	ND	18,000
PCB 1221	ND	18,000
PCB 1232	ND	18,000
PCB 1242	120,000	18,000
PCB 1248	ND	18,000
PCB 1254	ND	18,000
PCB 1260	81,000	18,000
% Surrogate Recovery:	%1	
Date Extracted:	10/19/94	
Date Analyzed:	10/23/94	

¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate = DCB (decachlorobiphenyl)

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-12, Motor Pile South of S-11 - EW Plot

TCT ID:	43012	
Parameter:		POL
PCB 1016	ND	19,000
PCB 1221	ND	19,000
PCB 1232	ND	19,000
PCB 1242	87,000	19,000
PCB 1248	ND	19,000
PCB 1254	ND	19,000
PCB 1260	37,000	19,000
% Surrogate Recovery:	%1	
Date Extracted:	10/19/94	
Date Analyzed:	10/23/94	

¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate = DCB (decachlorobiphenyl)

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-13, Large Ferrous Metal Pile - EW Plot

TCT ID:	43013	
Parameter:		<u>POL</u>
PCB 1016	ND	1,500
PCB 1221	ND	1,500
PCB 1232	ND	1,500
PCB 1242	5,300	1,500
PCB 1248	ND	1,500
PCB 1254	ND	1,500
PCB 1260	1,800	1,500
% Surrogate Recovery:	%1	
Date Extracted:	10/19/94	
Date Extracted:	10/17/74	
Date Analyzed:	10/23/94	

¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate = DCB (decachlorobiphenyl)

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-14, Copper Fines Pile, East of S-13 - EW Plot

TCT ID:	43014	·
Parameter:		PQL
PCB 1016	ND	17,000
PCB 1221	ND	17,000
PCB 1232	ND	17,000
PCB 1242	170,000	17,000
PCB 1248	ND	17,000
PCB 1254	ND	17,000
PCB 1260	ND	17,000
% Surrogate Recovery:	%1	
Date Extracted:	10/19/94	
Date Analyzed:	10/23/94	
Date Mialyzeu.	10/43/74	

¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

Surrogate = DCB (decachlorobiphenyl)

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in µg/Kg which is equal to parts-per-billion)

Client ID:

S-16, Process Area by Conveyor - East Plot

TCT ID:	43016	
Parameter:		POL
PCB 1016	ND	32,000
PCB 1221	ND	32,000
PCB 1232	ND ·	32,000
PCB 1242	270,000	32,000
PCB 1248	ND	32,000
PCB 1254	ND	32,000
PCB 1260	ND	32,000
% Surrogate Recovery:	% ¹	
Data E to and	10/10/04	
Date Extracted:	10/19/94	
Date Analyzed:	10/23/94	

¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit ND = Not Detected

Surrogate = DCB (decachlorobiphenyl)

POLYCHLORINATED BIPHENYL RESULTS EPA METHOD 8080

(All values are in $\mu g/Kg$ which is equal to parts-per-billion)

Client ID:

S-15, Sorted Steel Pile, East of S-14 - EW Plot

TCT ID:	43015	
Parameter:		POL
PCB 1016	ND	8,200
PCB 1221	ND	8,200
PCB 1232	ND	8,200
PCB 1242	33,000	8,200
PCB 1248	ND	8,200
PCB 1254	ND .	8,200
PCB 1260	ND	8,200
% Surrogate Recovery:	%1	
Date Futureted.	10/10/04	
Date Extracted:	10/19/94	
Date Analyzed:	10/23/94	

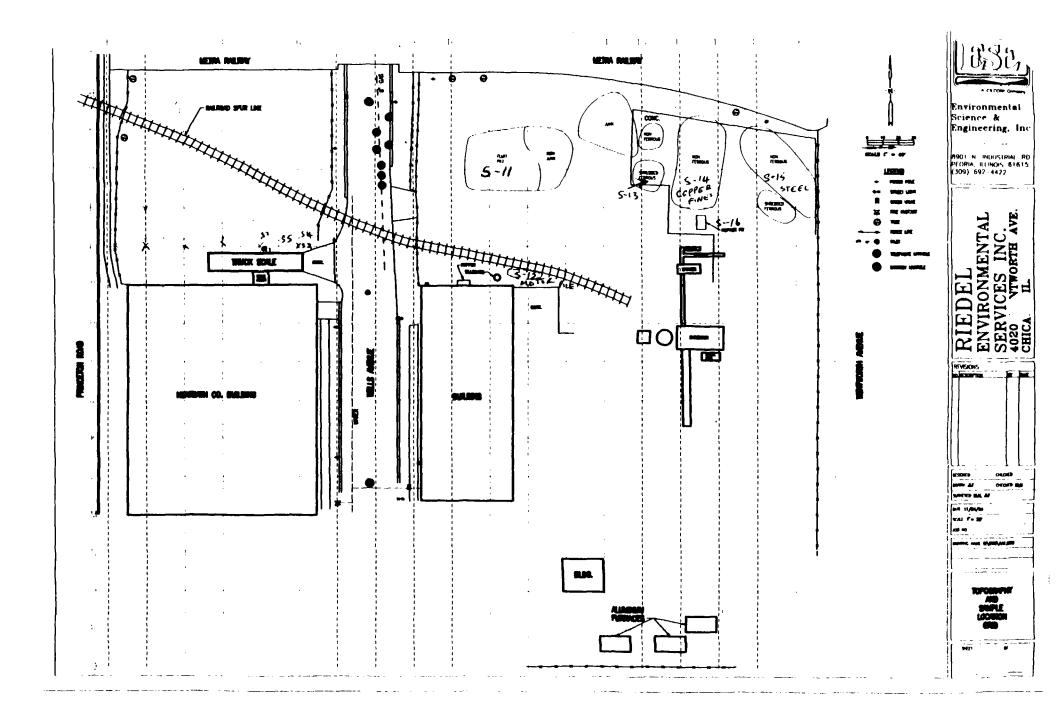
¹ Surrogate diluted out.

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

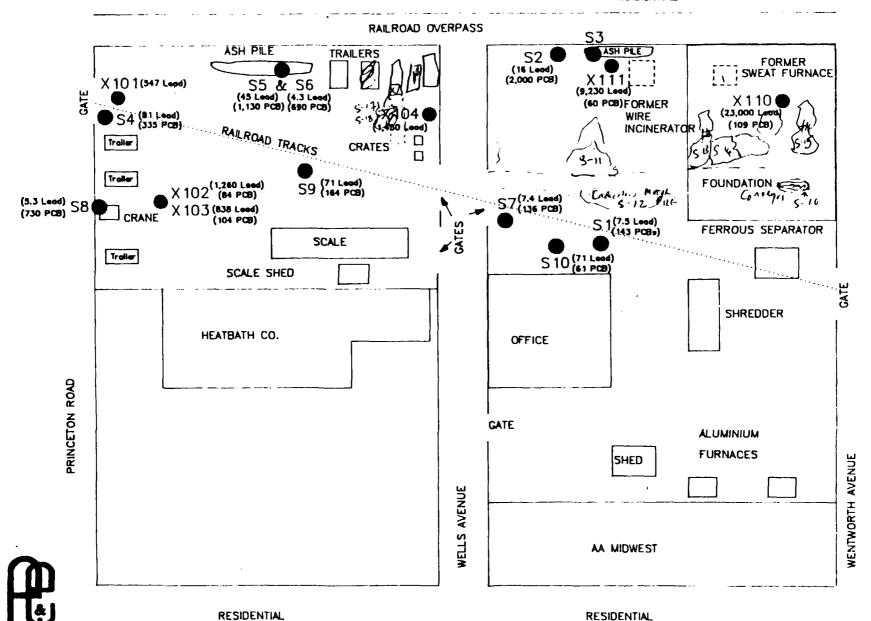
Surrogate = DCB (decachlorobiphenyl)



(

RED = TAT COLLECTED SAMPLES (Load = TCLP in mg/L; PCB = PPM)

BLUE = STATE COLLECTED SAMPLES (Lead = Totals in PPM; PCB = PPM)



RED = TAT COLLECTED SAMPLES (Lead = TCLP in mg/L; PCB = PPM)

BLUE = STATE COLLECTED SAMPLES (Load = Totals in PPM; PCB = PPM)

